

Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN060824\  
 Data File : BN031856.D  
 Acq On : 08 Jun 2024 13:14  
 Operator : MA/JU  
 Sample : P2634-15  
 Misc :  
 ALS Vial : 7 Sample Multiplier: 1

Instrument :  
 BNA\_N  
 ClientSampleId :  
 BHBX9

Integration Parameters: LSCINT.P

Integrator: RTE

Smoothing : OFF

Filtering: 5

Sampling : 1

Min Area: 0.1 % of largest Peak

Start Thrs: 0.2

Max Peaks: 100

Stop Thrs : 0

Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >

Peak separation: 5

Method : Z:\svoasrv\HPCHEM1\BNA\_N\Methods\SFAM-EPA-BN052924.MA.M  
 Title : SVOA CALIBRATION

Signal : TIC: BN031856.D\data.ms

| peak<br># | R.T.<br>min | first<br>scan | max<br>scan | last<br>scan | PK<br>TY | peak<br>height | corr.<br>area | corr.<br>% max. | % of<br>total |
|-----------|-------------|---------------|-------------|--------------|----------|----------------|---------------|-----------------|---------------|
| 1         | 3.593       | 99            | 103         | 115          | rBV      | 374133         | 581103        | 7.45%           | 0.370%        |
| 2         | 3.693       | 115           | 120         | 128          | rVV      | 293187         | 394278        | 5.06%           | 0.251%        |
| 3         | 4.893       | 319           | 324         | 330          | rVV2     | 114427         | 192743        | 2.47%           | 0.123%        |
| 4         | 5.675       | 452           | 457         | 465          | rVV      | 375922         | 669865        | 8.59%           | 0.427%        |
| 5         | 6.846       | 645           | 656         | 662          | rVV      | 1440040        | 2547356       | 32.67%          | 1.624%        |
| 6         | 7.010       | 669           | 684         | 696          | rVV      | 1154890        | 1876634       | 24.07%          | 1.196%        |
| 7         | 7.205       | 710           | 717         | 725          | rVV      | 1479833        | 2400168       | 30.79%          | 1.530%        |
| 8         | 7.281       | 725           | 730         | 734          | rVV2     | 94819          | 179620        | 2.30%           | 0.115%        |
| 9         | 7.669       | 787           | 796         | 803          | rVV      | 1433247        | 2381800       | 30.55%          | 1.518%        |
| 10        | 8.375       | 910           | 916         | 924          | rBV      | 1478227        | 2489351       | 31.93%          | 1.587%        |
| 11        | 8.828       | 985           | 993         | 999          | rBV      | 1369908        | 2304490       | 29.56%          | 1.469%        |
| 12        | 9.393       | 1083          | 1089        | 1102         | rVB      | 174154         | 300555        | 3.86%           | 0.192%        |
| 13        | 9.540       | 1106          | 1114        | 1126         | rBV      | 1121971        | 1996225       | 25.61%          | 1.273%        |
| 14        | 10.081      | 1199          | 1206        | 1215         | rVV      | 1735342        | 3043417       | 39.04%          | 1.940%        |
| 15        | 10.451      | 1259          | 1269        | 1275         | rBV      | 2000583        | 3554617       | 45.59%          | 2.266%        |
| 16        | 10.504      | 1275          | 1278        | 1283         | rVV      | 124861         | 187803        | 2.41%           | 0.120%        |
| 17        | 10.598      | 1284          | 1294        | 1315         | rVV      | 487648         | 1059111       | 13.59%          | 0.675%        |
| 18        | 11.198      | 1390          | 1396        | 1402         | rBV      | 269651         | 491336        | 6.30%           | 0.313%        |
| 19        | 12.122      | 1547          | 1553        | 1562         | rVB      | 176210         | 314035        | 4.03%           | 0.200%        |
| 20        | 12.340      | 1584          | 1590        | 1599         | rBV      | 120823         | 222837        | 2.86%           | 0.142%        |
| 21        | 13.134      | 1720          | 1725        | 1733         | rVB2     | 185813         | 313331        | 4.02%           | 0.200%        |
| 22        | 13.634      | 1805          | 1810        | 1815         | rVV      | 224815         | 371802        | 4.77%           | 0.237%        |
| 23        | 13.728      | 1821          | 1826        | 1832         | rVB      | 2920217        | 4199151       | 53.86%          | 2.677%        |
| 24        | 14.004      | 1867          | 1873        | 1887         | rVB      | 3611866        | 6280013       | 80.55%          | 4.003%        |
| 25        | 14.216      | 1905          | 1909        | 1916         | rVB      | 161175         | 211130        | 2.71%           | 0.135%        |
| 26        | 14.316      | 1916          | 1926        | 1933         | rBV      | 3013525        | 5018291       | 64.37%          | 3.199%        |
| 27        | 14.534      | 1956          | 1963        | 1970         | rBV      | 1351542        | 2217133       | 28.44%          | 1.413%        |
| 28        | 14.681      | 1983          | 1988        | 1991         | rVV2     | 261269         | 396012        | 5.08%           | 0.252%        |
| 29        | 14.716      | 1991          | 1994        | 2002         | rVV      | 268794         | 433478        | 5.56%           | 0.276%        |
| 30        | 14.792      | 2002          | 2007        | 2011         | rVB      | 171294         | 236497        | 3.03%           | 0.151%        |
| 31        | 14.986      | 2035          | 2040        | 2044         | rVB      | 181105         | 249139        | 3.20%           | 0.159%        |
| 32        | 15.104      | 2054          | 2060        | 2064         | rBV3     | 116145         | 191129        | 2.45%           | 0.122%        |
| 33        | 15.310      | 2088          | 2095        | 2101         | rVV      | 4789268        | 7144214       | 91.64%          | 4.554%        |
| 34        | 15.363      | 2101          | 2104        | 2107         | rVV2     | 139712         | 200531        | 2.57%           | 0.128%        |
| 35        | 15.398      | 2107          | 2110        | 2113         | rVV      | 159582         | 228032        | 2.92%           | 0.145%        |
| 36        | 15.439      | 2113          | 2117        | 2122         | rVV      | 744259         | 1025683       | 13.16%          | 0.654%        |

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 Misc :  
 ALS Vial : 7 Sample Multiplier: 1

Instrument :  
 BNA\_N  
 ClientSampleId :  
 BHBX9

Integration Parameters: LSCINT.P

Integrator: RTE

Smoothing : OFF

Filtering: 5

Sampling : 1

Min Area: 0.1 % of largest Peak

Start Thrs: 0.2

Max Peaks: 100

Stop Thrs : 0

Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >

Peak separation: 5

Method : Z:\svoasrv\HPCHEM1\BNA\_N\Methods\SFAM-EPA-BN052924.MA.M  
 Title : SVOA CALIBRATION

|    |        |      |      |      |      |         |         |        |        |
|----|--------|------|------|------|------|---------|---------|--------|--------|
| 37 | 15.657 | 2149 | 2154 | 2161 | rVV2 | 146784  | 256760  | 3.29%  | 0.164% |
| 38 | 15.810 | 2172 | 2180 | 2186 | rVB3 | 71639   | 191732  | 2.46%  | 0.122% |
| 39 | 16.034 | 2212 | 2218 | 2226 | rVB5 | 268486  | 659655  | 8.46%  | 0.421% |
| 40 | 16.369 | 2267 | 2275 | 2279 | rBV3 | 134195  | 339114  | 4.35%  | 0.216% |
| 41 | 16.416 | 2279 | 2283 | 2289 | rVV  | 342796  | 497125  | 6.38%  | 0.317% |
| 42 | 16.669 | 2322 | 2326 | 2331 | rVB  | 714590  | 963518  | 12.36% | 0.614% |
| 43 | 16.722 | 2331 | 2335 | 2342 | rVB2 | 413744  | 659185  | 8.46%  | 0.420% |
| 44 | 16.822 | 2342 | 2352 | 2355 | rBV4 | 142369  | 381661  | 4.90%  | 0.243% |
| 45 | 16.881 | 2358 | 2362 | 2370 | rVB  | 257527  | 429667  | 5.51%  | 0.274% |
| 46 | 17.063 | 2384 | 2393 | 2397 | rBV  | 3573511 | 5357363 | 68.72% | 3.415% |
| 47 | 17.104 | 2397 | 2400 | 2405 | rVV  | 2902585 | 3896198 | 49.98% | 2.484% |
| 48 | 17.163 | 2405 | 2410 | 2414 | rVV2 | 4829005 | 7018675 | 90.03% | 4.474% |
| 49 | 17.198 | 2414 | 2416 | 2420 | rVB  | 501239  | 545348  | 7.00%  | 0.348% |
| 50 | 17.257 | 2420 | 2426 | 2430 | rBV3 | 216764  | 402754  | 5.17%  | 0.257% |
| 51 | 17.345 | 2436 | 2441 | 2443 | rBV2 | 145095  | 239636  | 3.07%  | 0.153% |
| 52 | 17.375 | 2443 | 2446 | 2451 | rVB3 | 321391  | 481105  | 6.17%  | 0.307% |
| 53 | 17.469 | 2454 | 2462 | 2467 | rBV  | 271718  | 532807  | 6.83%  | 0.340% |
| 54 | 17.645 | 2488 | 2492 | 2500 | rVB2 | 414028  | 642225  | 8.24%  | 0.409% |
| 55 | 17.745 | 2505 | 2509 | 2512 | rBV2 | 150238  | 212415  | 2.72%  | 0.135% |
| 56 | 17.786 | 2513 | 2516 | 2521 | rVB  | 113601  | 182422  | 2.34%  | 0.116% |
| 57 | 17.857 | 2524 | 2528 | 2532 | rVB3 | 124703  | 180517  | 2.32%  | 0.115% |
| 58 | 17.939 | 2537 | 2542 | 2544 | rBV  | 925279  | 1251600 | 16.05% | 0.798% |
| 59 | 17.969 | 2544 | 2547 | 2548 | rBV  | 1174626 | 1189821 | 15.26% | 0.758% |
| 60 | 18.128 | 2569 | 2574 | 2578 | rVV2 | 903460  | 1580263 | 20.27% | 1.007% |
| 61 | 18.169 | 2578 | 2581 | 2588 | rVB  | 546691  | 775932  | 9.95%  | 0.495% |
| 62 | 18.257 | 2592 | 2596 | 2599 | rBV3 | 301359  | 450621  | 5.78%  | 0.287% |
| 63 | 18.292 | 2600 | 2602 | 2606 | rVV2 | 205039  | 287456  | 3.69%  | 0.183% |
| 64 | 18.333 | 2606 | 2609 | 2614 | rVB3 | 249737  | 366278  | 4.70%  | 0.233% |
| 65 | 18.445 | 2623 | 2628 | 2632 | rBV  | 676276  | 988100  | 12.67% | 0.630% |
| 66 | 18.486 | 2632 | 2635 | 2641 | rVB2 | 512106  | 715658  | 9.18%  | 0.456% |
| 67 | 18.692 | 2666 | 2670 | 2674 | rVV  | 338376  | 468912  | 6.01%  | 0.299% |
| 68 | 18.739 | 2674 | 2678 | 2682 | rVV  | 452867  | 646847  | 8.30%  | 0.412% |
| 69 | 18.780 | 2682 | 2685 | 2689 | rVB  | 335305  | 426839  | 5.48%  | 0.272% |
| 70 | 18.863 | 2694 | 2699 | 2703 | rBV  | 784080  | 1192939 | 15.30% | 0.760% |
| 71 | 18.957 | 2712 | 2715 | 2718 | rVB  | 253903  | 289925  | 3.72%  | 0.185% |
| 72 | 19.004 | 2719 | 2723 | 2729 | rBV2 | 452804  | 860082  | 11.03% | 0.548% |
| 73 | 19.127 | 2738 | 2744 | 2751 | rBV  | 5608855 | 7756455 | 99.49% | 4.944% |
| 74 | 19.192 | 2752 | 2755 | 2758 | rBV  | 318694  | 372104  | 4.77%  | 0.237% |
| 75 | 19.251 | 2762 | 2765 | 2768 | rBV3 | 332882  | 357785  | 4.59%  | 0.228% |
| 76 | 19.463 | 2796 | 2801 | 2804 | rBV  | 4841632 | 7664291 | 98.31% | 4.886% |

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 Sample : P2634-15  
 Misc :  
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Instrument :  
 BNA\_N  
 ClientSampleId :  
 BHBX9

Integration Parameters: LSCINT.P

Integrator: RTE

Smoothing : OFF

Filtering: 5

Sampling : 1

Min Area: 0.1 % of largest Peak

Start Thrs: 0.2

Max Peaks: 100

Stop Thrs : 0

Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >

Peak separation: 5

Method : Z:\svoasrv\HPCHEM1\BNA\_N\Methods\SFAM-EPA-BN052924.MA.M  
 Title : SVOA CALIBRATION

|     |        |      |      |      |      |         |         |         |        |
|-----|--------|------|------|------|------|---------|---------|---------|--------|
| 77  | 19.492 | 2804 | 2806 | 2814 | rVV  | 4400469 | 5722505 | 73.40%  | 3.648% |
| 78  | 19.569 | 2815 | 2819 | 2824 | rVB2 | 422461  | 649258  | 8.33%   | 0.414% |
| 79  | 19.733 | 2844 | 2847 | 2852 | rVB3 | 263937  | 365291  | 4.69%   | 0.233% |
| 80  | 19.816 | 2856 | 2861 | 2872 | rBV4 | 489203  | 1264844 | 16.22%  | 0.806% |
| 81  | 19.951 | 2880 | 2884 | 2887 | rBV2 | 367676  | 663528  | 8.51%   | 0.423% |
| 82  | 19.992 | 2887 | 2891 | 2894 | rVB  | 601899  | 849534  | 10.90%  | 0.542% |
| 83  | 20.092 | 2904 | 2908 | 2911 | rBV  | 487692  | 588531  | 7.55%   | 0.375% |
| 84  | 20.151 | 2914 | 2918 | 2921 | rVB2 | 511201  | 689247  | 8.84%   | 0.439% |
| 85  | 20.286 | 2938 | 2941 | 2945 | rBV  | 394605  | 520776  | 6.68%   | 0.332% |
| 86  | 20.333 | 2947 | 2949 | 2954 | rVB  | 339917  | 379415  | 4.87%   | 0.242% |
| 87  | 20.739 | 3015 | 3018 | 3024 | rBV  | 608693  | 945004  | 12.12%  | 0.602% |
| 88  | 20.969 | 3052 | 3057 | 3059 | rBV2 | 1767336 | 2714291 | 34.82%  | 1.730% |
| 89  | 21.210 | 3096 | 3098 | 3103 | rVB  | 507894  | 597028  | 7.66%   | 0.381% |
| 90  | 21.298 | 3106 | 3113 | 3118 | rVV4 | 3360030 | 7796100 | 100.00% | 4.970% |
| 91  | 21.345 | 3118 | 3121 | 3133 | rVB  | 2428360 | 3776417 | 48.44%  | 2.407% |
| 92  | 21.504 | 3142 | 3148 | 3153 | rBV3 | 1581812 | 2737541 | 35.11%  | 1.745% |
| 93  | 21.569 | 3155 | 3159 | 3164 | rVV  | 2065584 | 3183335 | 40.83%  | 2.029% |
| 94  | 21.621 | 3165 | 3168 | 3171 | rVV  | 680845  | 972573  | 12.48%  | 0.620% |
| 95  | 21.663 | 3172 | 3175 | 3184 | rVB3 | 777647  | 1252400 | 16.06%  | 0.798% |
| 96  | 22.051 | 3237 | 3241 | 3252 | rVB6 | 530864  | 1166932 | 14.97%  | 0.744% |
| 97  | 22.668 | 3342 | 3346 | 3357 | rVB4 | 877453  | 1754623 | 22.51%  | 1.119% |
| 98  | 23.321 | 3451 | 3457 | 3463 | rBV  | 1555262 | 4132221 | 53.00%  | 2.634% |
| 99  | 24.051 | 3575 | 3581 | 3586 | rBV2 | 1372316 | 2722972 | 34.93%  | 1.736% |
| 100 | 24.251 | 3608 | 3615 | 3623 | rBV2 | 1589608 | 3813226 | 48.91%  | 2.431% |

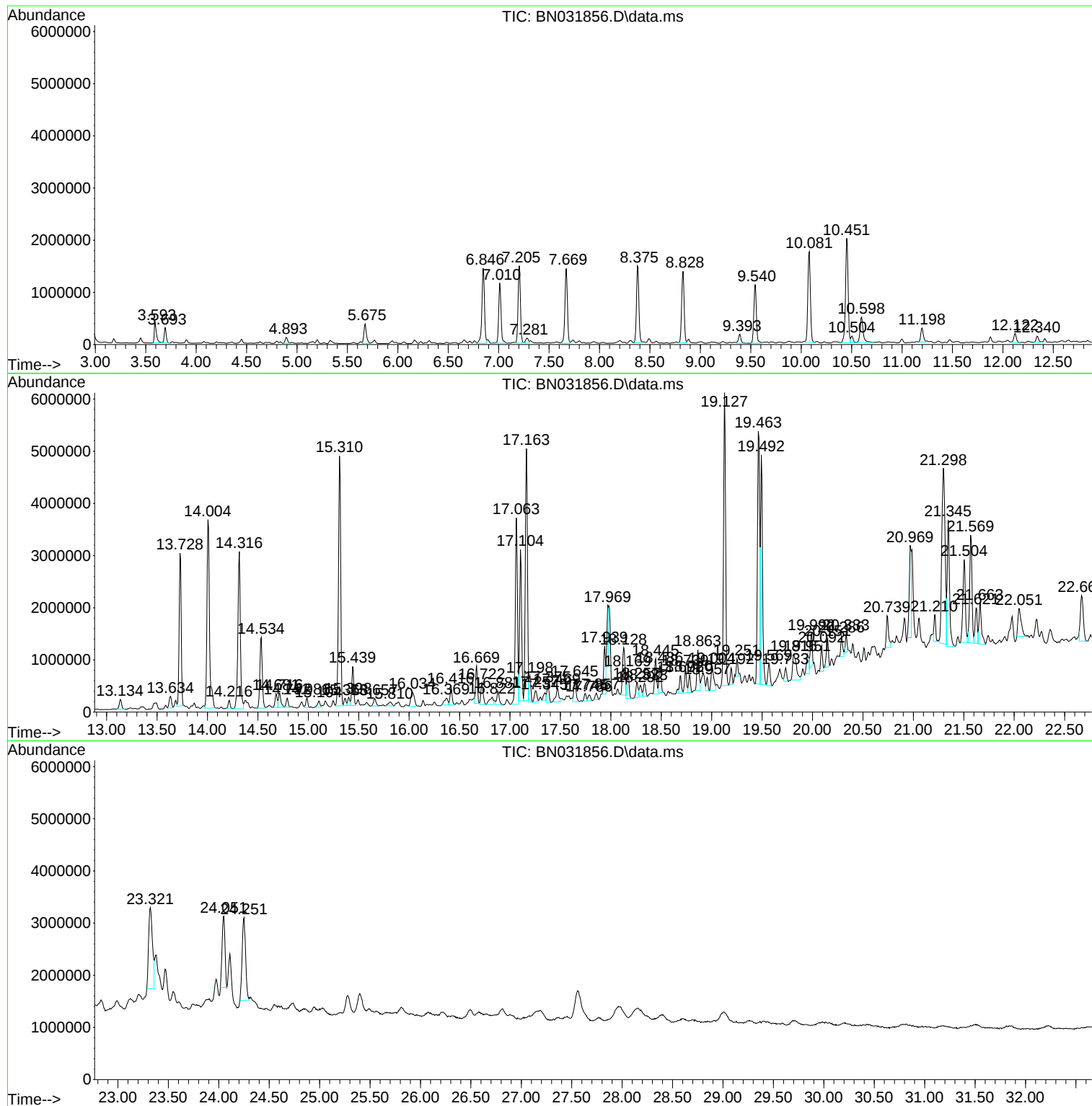
Sum of corrected areas: 156872292

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Sample : P2634-15  
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ALS Vial : 7 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
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Quant Method : Z:\svoasrv\HPCHEM1\BNA\_N\Methods\SFAM-EPA-BN052924.MA.M  
Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P



Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN060824\  
Data File : BN031856.D  
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Quant Method : Z:\svoasrv\HPCHEM1\BNA\_N\Methods\SFAM-EPA-BN052924.MA.M  
Quant Title : SVOA CALIBRATION

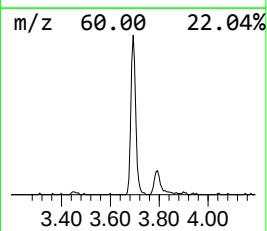
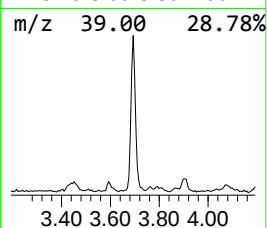
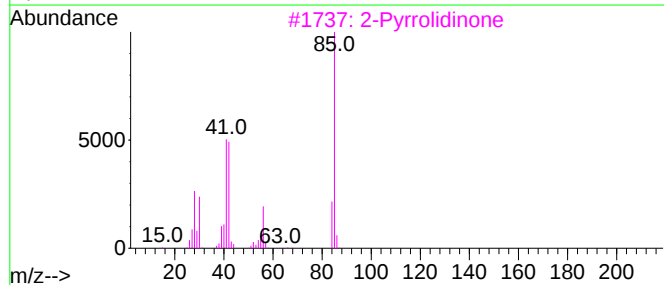
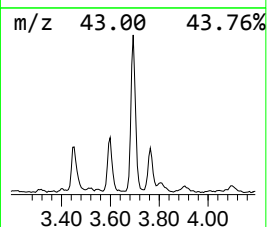
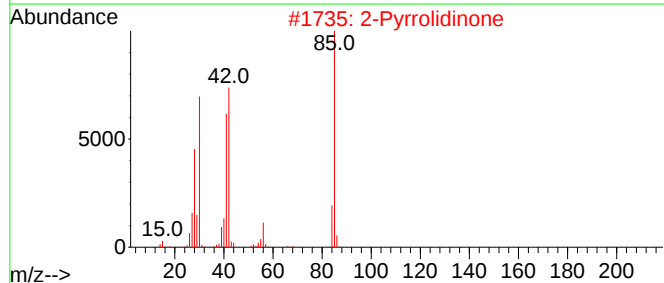
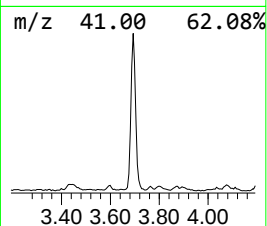
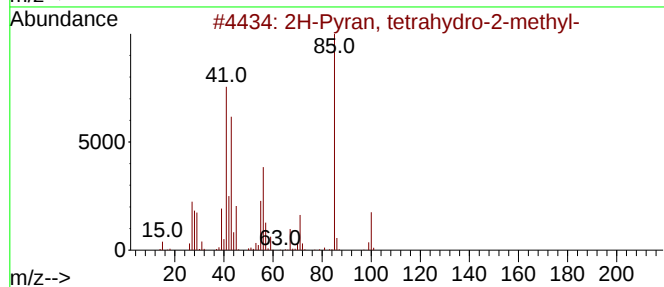
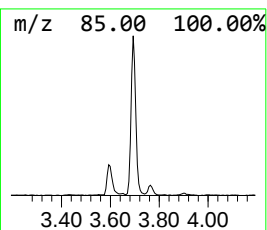
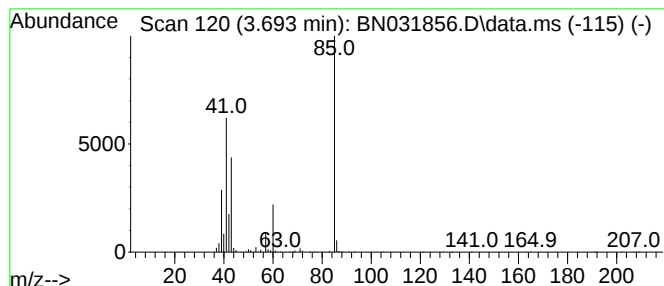
TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 1 unknown-01 Concentration Rank 12

| R.T.  | EstConc    | Area   | Relative to ISTD       | R.T.  |
|-------|------------|--------|------------------------|-------|
| 3.693 | 3.31 ng/ul | 394278 | 1,4-Dichlorobenzene-d4 | 7.669 |

| Hit# | of                             | 5 | Tentative ID | MW  | MolForm | CAS#        | Qual |
|------|--------------------------------|---|--------------|-----|---------|-------------|------|
| 1    | 2H-Pyran, tetrahydro-2-methyl- |   |              | 100 | C6H12O  | 010141-72-7 | 36   |
| 2    | 2-Pyrrolidinone                |   |              | 85  | C4H7NO  | 000616-45-5 | 33   |
| 3    | 2-Pyrrolidinone                |   |              | 85  | C4H7NO  | 000616-45-5 | 33   |
| 4    | 2-Pyrrolidinone                |   |              | 85  | C4H7NO  | 000616-45-5 | 9    |
| 5    | 2H-Pyran, tetrahydro-2-methyl- |   |              | 100 | C6H12O  | 010141-72-7 | 9    |



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ALS Vial : 7 Sample Multiplier: 1

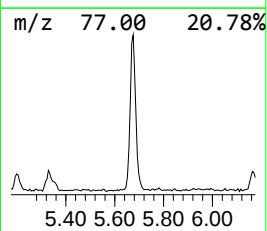
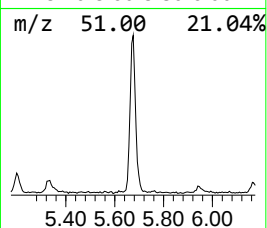
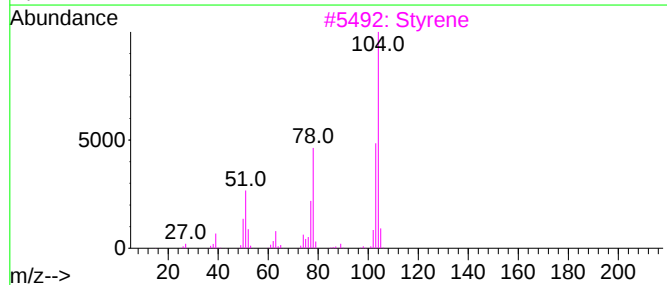
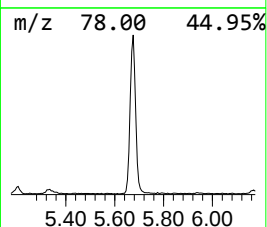
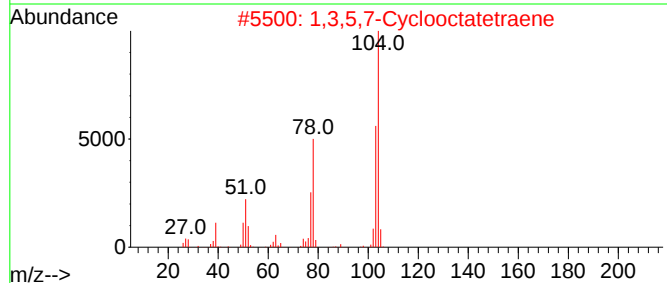
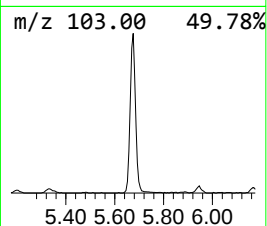
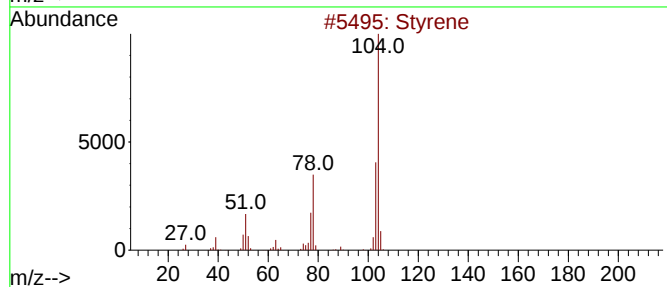
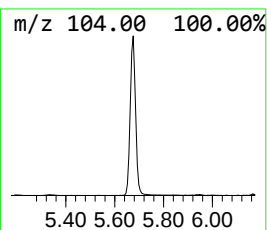
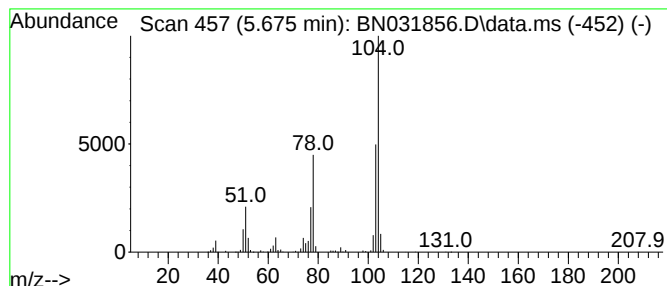
Instrument :  
BNA\_N  
ClientSampleId :  
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Quant Method : Z:\svoasrv\HPCHEM1\BNA\_N\Methods\SFAM-EPA-BN052924.MA.M  
Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 2 Styrene Concentration Rank 5

| R.T.      | EstConc                         | Area   | Relative to ISTD       |         |             | R.T.  |
|-----------|---------------------------------|--------|------------------------|---------|-------------|-------|
| 5.675     | 5.62 ng/ul                      | 669865 | 1,4-Dichlorobenzene-d4 |         |             | 7.669 |
| Hit# of 5 | Tentative ID                    |        | MW                     | MolForm | CAS#        | Qual  |
| 1         | Styrene                         |        | 104                    | C8H8    | 000100-42-5 | 97    |
| 2         | 1,3,5,7-Cyclooctatetraene       |        | 104                    | C8H8    | 000629-20-9 | 97    |
| 3         | Styrene                         |        | 104                    | C8H8    | 000100-42-5 | 96    |
| 4         | Bicyclo[4.2.0]octa-1,3,5-triene |        | 104                    | C8H8    | 000694-87-1 | 95    |
| 5         | 1,3,5,7-Cyclooctatetraene       |        | 104                    | C8H8    | 000629-20-9 | 95    |



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Data File : BN031856.D  
Acq On : 08 Jun 2024 13:14  
Operator : MA/JU  
Sample : P2634-15  
Misc :  
ALS Vial : 7 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
BHBX9

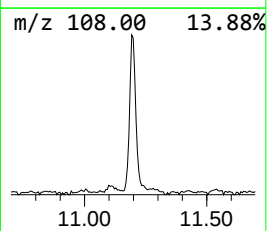
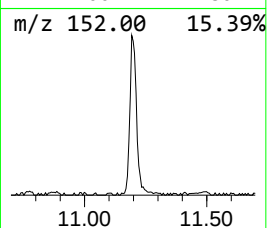
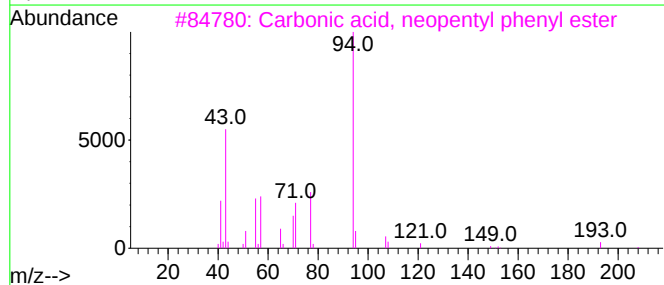
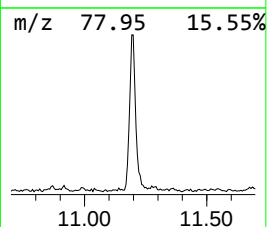
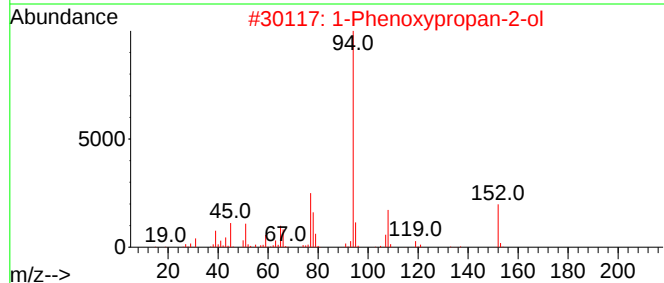
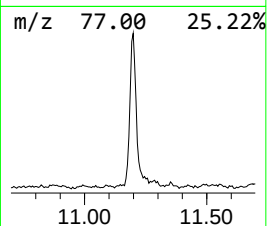
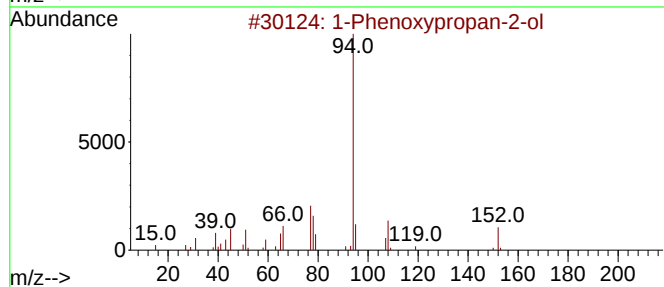
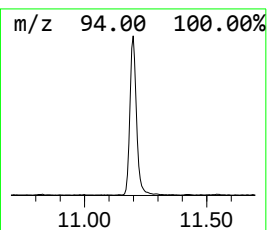
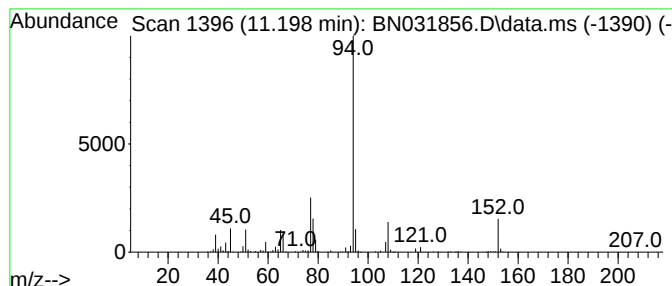
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Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 3 1-Phenoxypropan-2-ol Concentration Rank 18

| R.T.   | EstConc    | Area   | Relative to ISTD | R.T.   |
|--------|------------|--------|------------------|--------|
| 11.198 | 2.76 ng/ul | 491336 | Naphthalene-d8   | 10.451 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm  | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|----------|-------------|------|
| 1    |    |   | 1-Phenoxypropan-2-ol                | 152 | C9H12O2  | 000770-35-4 | 94   |
| 2    |    |   | 1-Phenoxypropan-2-ol                | 152 | C9H12O2  | 000770-35-4 | 70   |
| 3    |    |   | Carbonic acid, neopentyl phenyl ... | 208 | C12H16O3 | 013183-19-2 | 64   |
| 4    |    |   | 1-Propanol, 2-phenoxy-              | 152 | C9H12O2  | 004169-04-4 | 64   |
| 5    |    |   | Ethanol, 2-phenoxy-                 | 138 | C8H10O2  | 000122-99-6 | 64   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN060824\  
Data File : BN031856.D  
Acq On : 08 Jun 2024 13:14  
Operator : MA/JU  
Sample : P2634-15  
Misc :  
ALS Vial : 7 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
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Quant Title : SVOA CALIBRATION

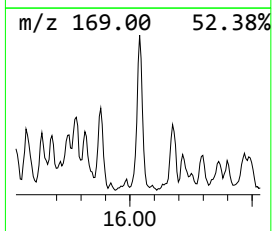
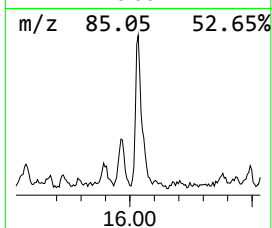
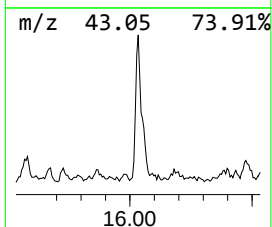
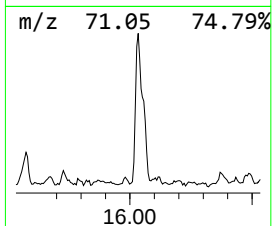
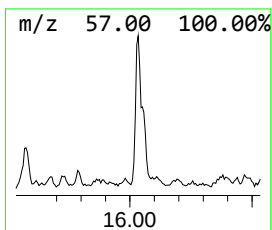
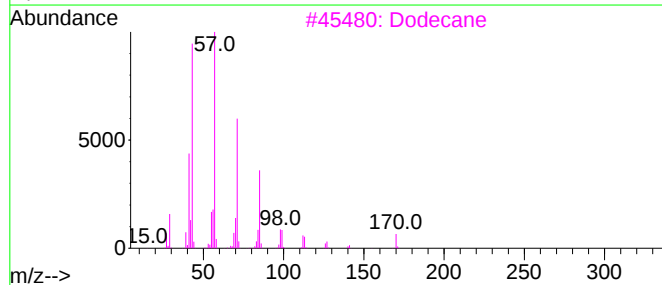
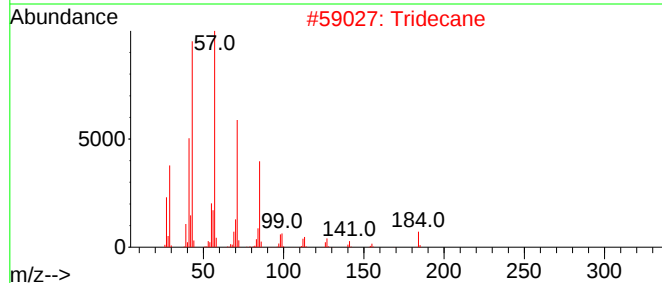
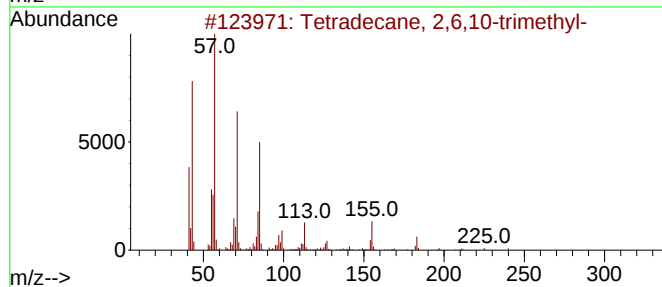
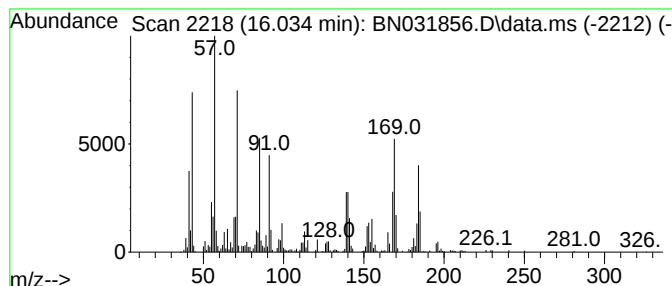
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TIC Integration Parameters: LSCINT.P

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Peak Number 4 (DEL) Alkane: Straight-Chai... Concentration Rank 21

| R.T.   | EstConc    | Area   | Relative to ISTD | R.T.   |
|--------|------------|--------|------------------|--------|
| 16.034 | 2.46 ng/ul | 659655 | Phenanthrene-d10 | 17.063 |

| Hit# | of | 5 | Tentative ID                   | MW  | MolForm | CAS#        | Qual |
|------|----|---|--------------------------------|-----|---------|-------------|------|
| 1    |    |   | Tetradecane, 2,6,10-trimethyl- | 240 | C17H36  | 014905-56-7 | 56   |
| 2    |    |   | Tridecane                      | 184 | C13H28  | 000629-50-5 | 55   |
| 3    |    |   | Dodecane                       | 170 | C12H26  | 000112-40-3 | 46   |
| 4    |    |   | Tridecane                      | 184 | C13H28  | 000629-50-5 | 46   |
| 5    |    |   | Dodecane                       | 170 | C12H26  | 000112-40-3 | 41   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN060824\  
Data File : BN031856.D  
Acq On : 08 Jun 2024 13:14  
Operator : MA/JU  
Sample : P2634-15  
Misc :  
ALS Vial : 7 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
BHBX9

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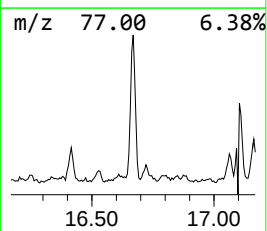
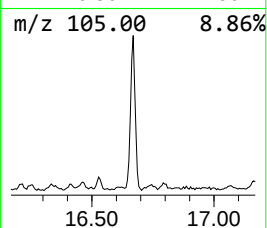
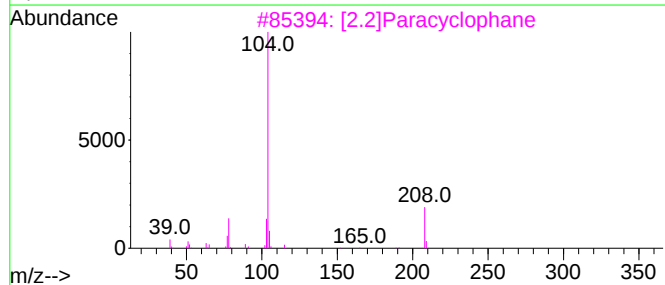
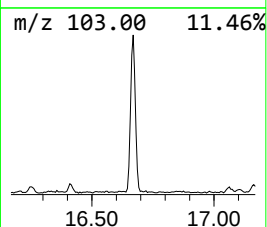
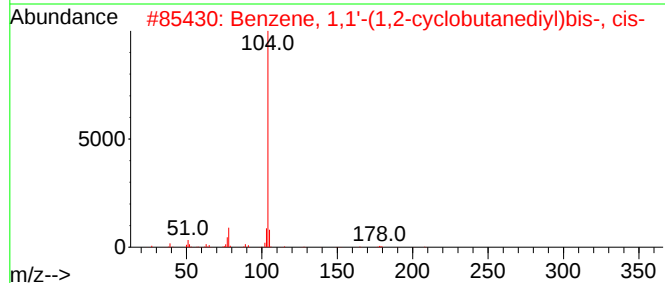
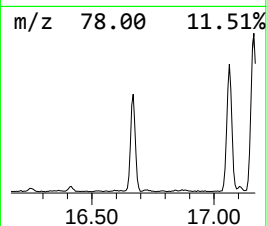
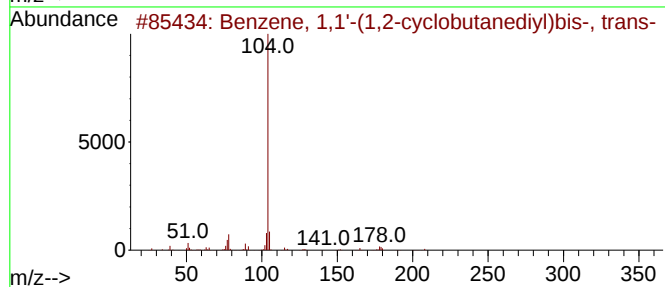
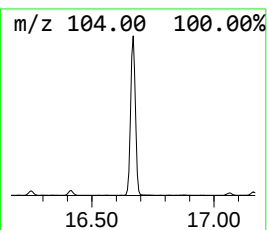
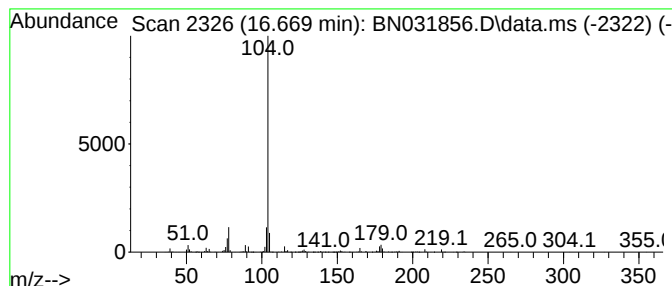
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TIC Integration Parameters: LSCINT.P

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Peak Number 5 Benzene, 1,1'-(1,2-cyclobut... Concentration Rank 11

| R.T.   | EstConc    | Area   | Relative to ISTD | R.T.   |
|--------|------------|--------|------------------|--------|
| 16.669 | 3.60 ng/ul | 963518 | Phenanthrene-d10 | 17.063 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|---------|-------------|------|
| 1    |    |   | Benzene, 1,1'-(1,2-cyclobutanedi... | 208 | C16H16  | 020071-09-4 | 87   |
| 2    |    |   | Benzene, 1,1'-(1,2-cyclobutanedi... | 208 | C16H16  | 007694-30-6 | 86   |
| 3    |    |   | [2.2]Paracyclophane                 | 208 | C16H16  | 001633-22-3 | 83   |
| 4    |    |   | Cyclobutane, 1,2-diphenyl-          | 208 | C16H16  | 003018-21-1 | 72   |
| 5    |    |   | Cyclobutane, 1,3-diphenyl-, trans-  | 208 | C16H16  | 025558-23-0 | 72   |



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Data File : BN031856.D  
Acq On : 08 Jun 2024 13:14  
Operator : MA/JU  
Sample : P2634-15  
Misc :  
ALS Vial : 7 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
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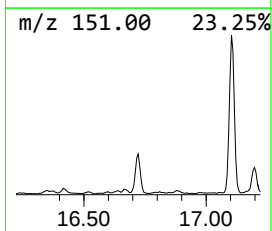
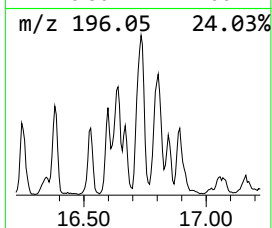
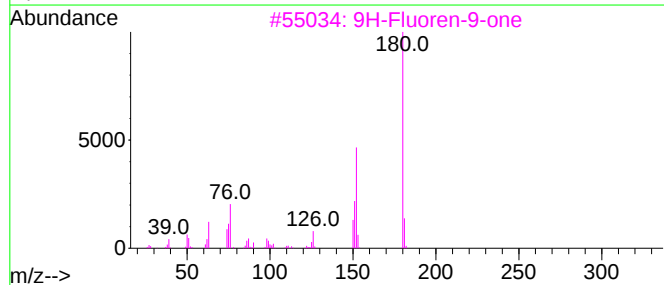
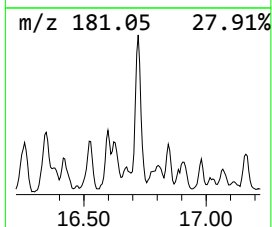
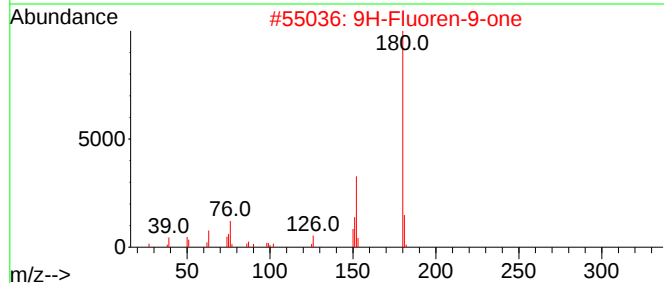
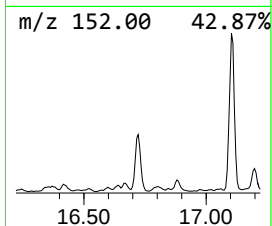
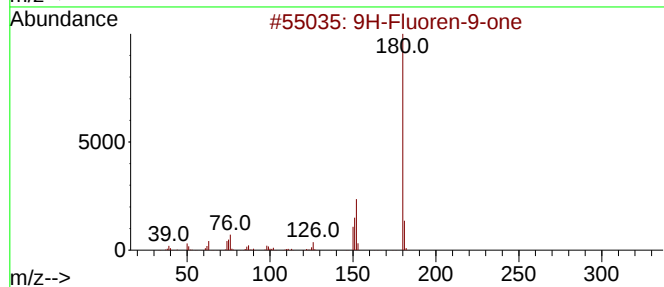
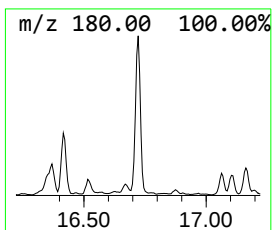
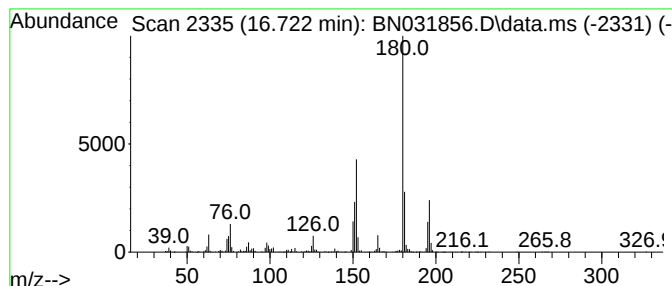
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TIC Integration Parameters: LSCINT.P

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Peak Number 6 9H-Fluoren-9-one Concentration Rank 22

| R.T.   | EstConc    | Area   | Relative to ISTD |  | R.T.   |
|--------|------------|--------|------------------|--|--------|
| 16.722 | 2.46 ng/ul | 659185 | Phenanthrene-d10 |  | 17.063 |

| Hit# | of 5 | Tentative ID           | MW  | MolForm | CAS#        | Qual |
|------|------|------------------------|-----|---------|-------------|------|
| 1    |      | 9H-Fluoren-9-one       | 180 | C13H8O  | 000486-25-9 | 93   |
| 2    |      | 9H-Fluoren-9-one       | 180 | C13H8O  | 000486-25-9 | 89   |
| 3    |      | 9H-Fluoren-9-one       | 180 | C13H8O  | 000486-25-9 | 76   |
| 4    |      | 9H-Fluoren-9-one       | 180 | C13H8O  | 000486-25-9 | 76   |
| 5    |      | 9,10-Phenanthrenedione | 208 | C14H8O2 | 000084-11-7 | 59   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN060824\  
Data File : BN031856.D  
Acq On : 08 Jun 2024 13:14  
Operator : MA/JU  
Sample : P2634-15  
Misc :  
ALS Vial : 7 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
BHBX9

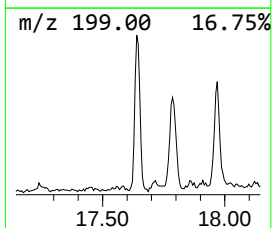
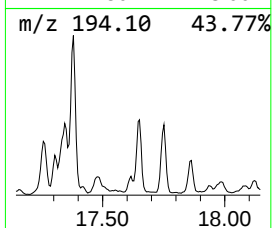
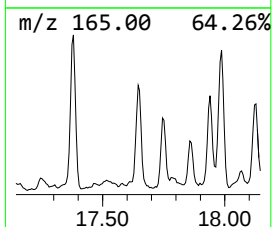
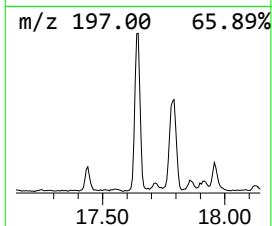
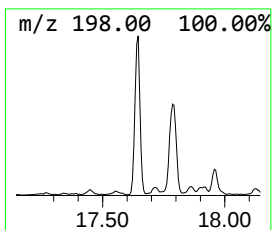
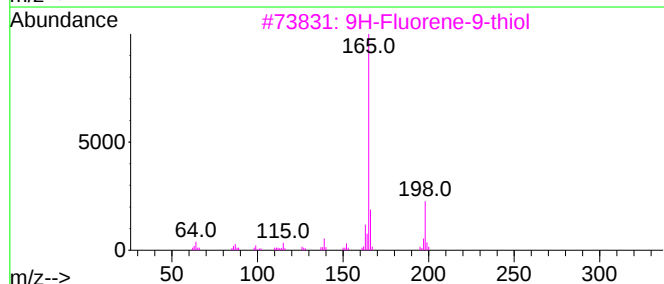
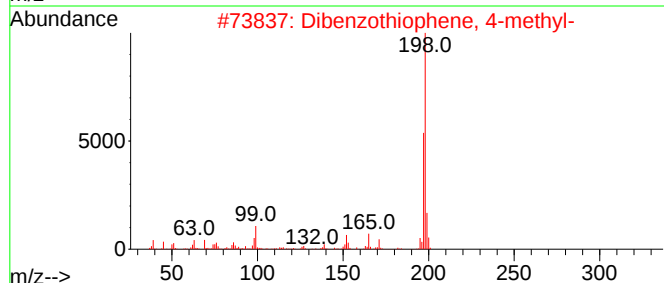
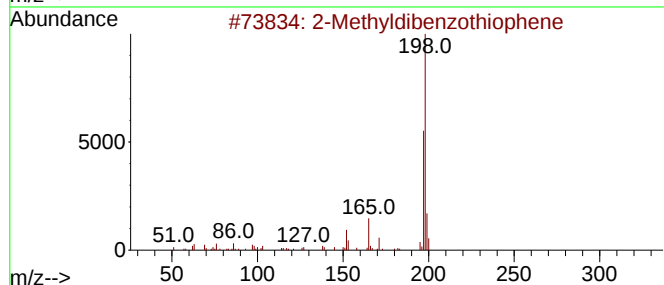
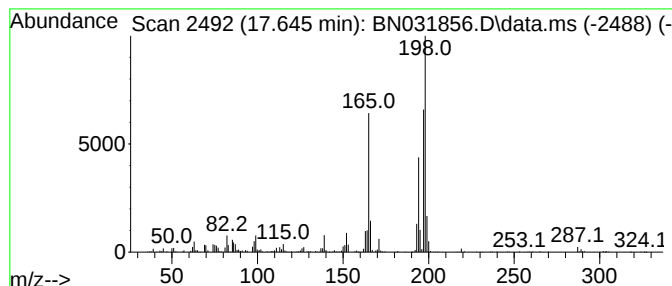
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Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 7 2-Methyldibenzothiophene Concentration Rank 25

| R.T.   | EstConc    | Area   | Relative to ISTD | R.T.   |
|--------|------------|--------|------------------|--------|
| 17.645 | 2.40 ng/ul | 642225 | Phenanthrene-d10 | 17.063 |

| Hit# | of | 5 | Tentative ID                | MW  | MolForm  | CAS#         | Qual |
|------|----|---|-----------------------------|-----|----------|--------------|------|
| 1    |    |   | 2-Methyldibenzothiophene    | 198 | C13H10S  | 1000383-55-1 | 68   |
| 2    |    |   | Dibenzothiophene, 4-methyl- | 198 | C13H10S  | 007372-88-5  | 55   |
| 3    |    |   | 9H-Fluorene-9-thiol         | 198 | C13H10S  | 019552-08-0  | 53   |
| 4    |    |   | Benzo[c]cinoline, 4-methyl- | 194 | C13H10N2 | 019174-78-8  | 51   |
| 5    |    |   | Dibenzothiophene, 3-methyl- | 198 | C13H10S  | 016587-52-3  | 49   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN060824\  
Data File : BN031856.D  
Acq On : 08 Jun 2024 13:14  
Operator : MA/JU  
Sample : P2634-15  
Misc :  
ALS Vial : 7 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
BHBX9

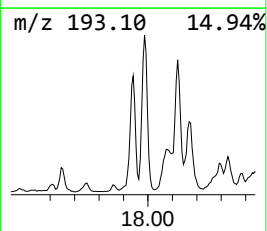
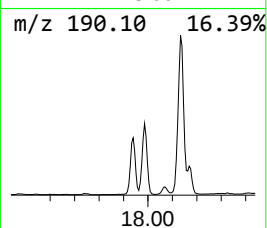
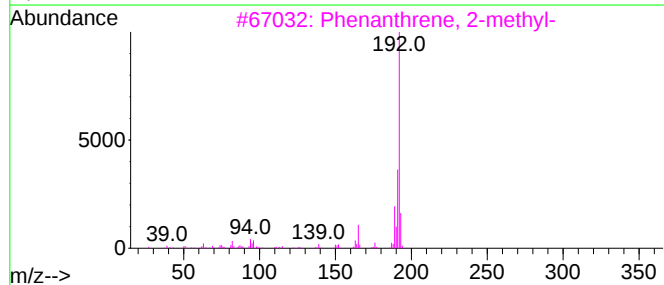
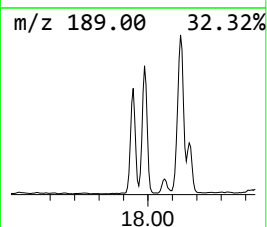
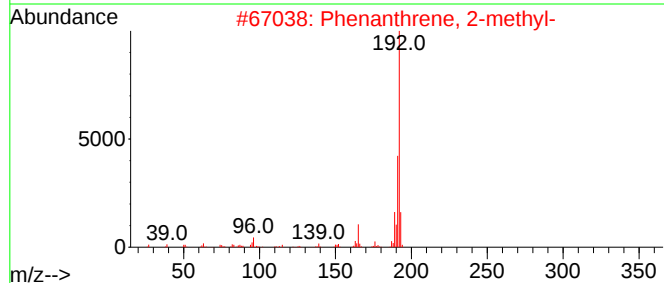
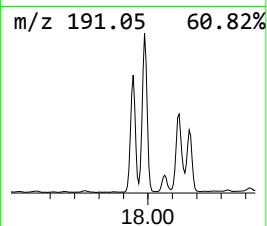
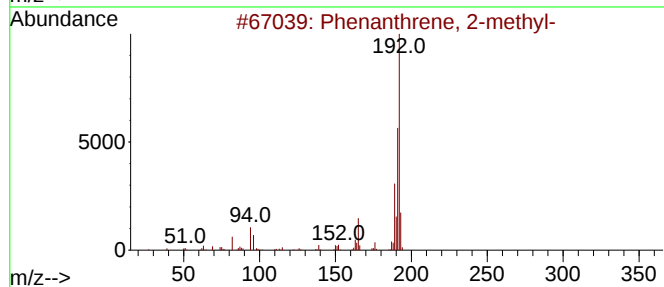
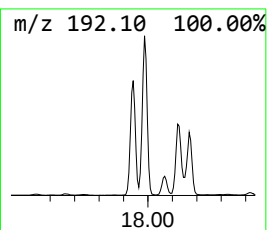
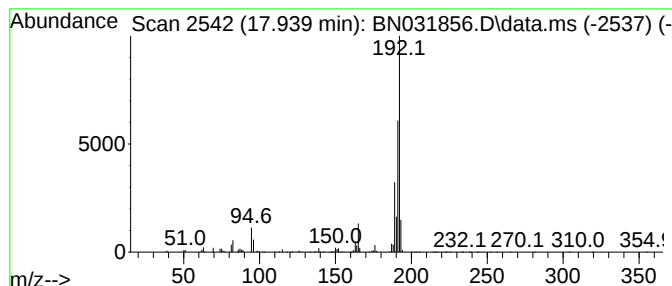
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Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 8 Phenanthrene, 2-methyl- Concentration Rank 6

| R.T.   | EstConc    | Area    | Relative to ISTD | R.T.   |
|--------|------------|---------|------------------|--------|
| 17.939 | 4.67 ng/ul | 1251600 | Phenanthrene-d10 | 17.063 |

| Hit# | of | 5 | Tentative ID            | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------|-----|---------|-------------|------|
| 1    |    |   | Phenanthrene, 2-methyl- | 192 | C15H12  | 002531-84-2 | 96   |
| 2    |    |   | Phenanthrene, 2-methyl- | 192 | C15H12  | 002531-84-2 | 96   |
| 3    |    |   | Phenanthrene, 2-methyl- | 192 | C15H12  | 002531-84-2 | 94   |
| 4    |    |   | Phenanthrene, 4-methyl- | 192 | C15H12  | 000832-64-4 | 94   |
| 5    |    |   | Phenanthrene, 1-methyl- | 192 | C15H12  | 000832-69-9 | 93   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN060824\  
Data File : BN031856.D  
Acq On : 08 Jun 2024 13:14  
Operator : MA/JU  
Sample : P2634-15  
Misc :  
ALS Vial : 7 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
BHBX9

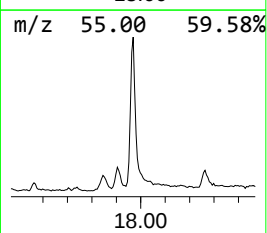
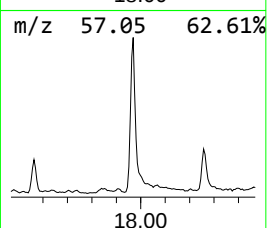
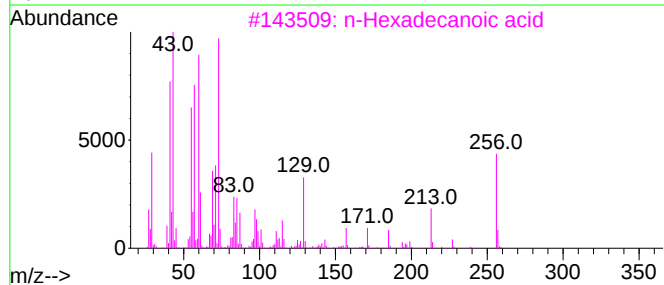
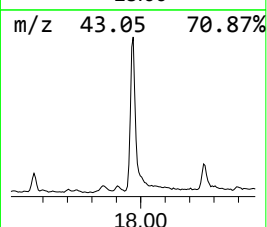
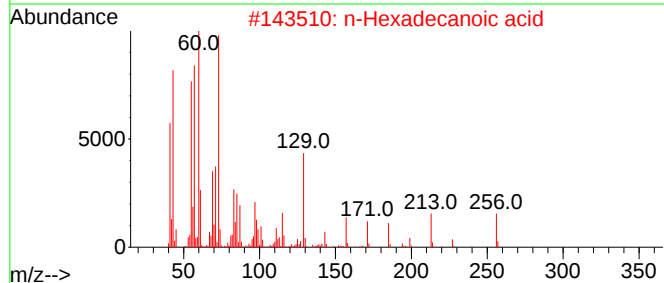
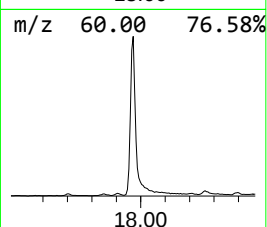
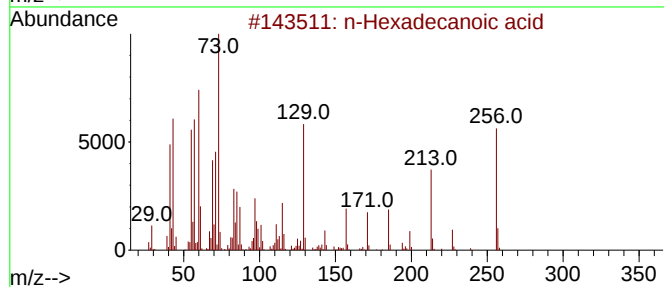
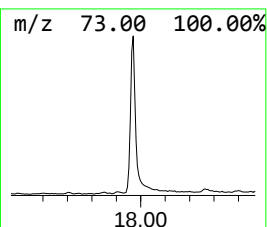
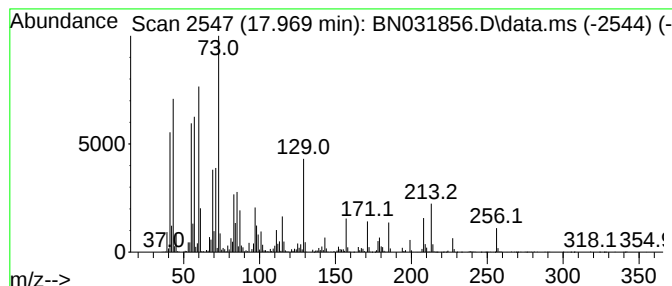
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Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 9 n-Hexadecanoic acid Concentration Rank 9

| R.T.   | EstConc    | Area    | Relative to ISTD | R.T.   |
|--------|------------|---------|------------------|--------|
| 17.969 | 4.44 ng/ul | 1189820 | Phenanthrene-d10 | 17.063 |

| Hit# | of | 5 | Tentative ID        | MW  | MolForm  | CAS#        | Qual |
|------|----|---|---------------------|-----|----------|-------------|------|
| 1    |    |   | n-Hexadecanoic acid | 256 | C16H32O2 | 000057-10-3 | 99   |
| 2    |    |   | n-Hexadecanoic acid | 256 | C16H32O2 | 000057-10-3 | 99   |
| 3    |    |   | n-Hexadecanoic acid | 256 | C16H32O2 | 000057-10-3 | 98   |
| 4    |    |   | n-Hexadecanoic acid | 256 | C16H32O2 | 000057-10-3 | 98   |
| 5    |    |   | Tetradecanoic acid  | 228 | C14H28O2 | 000544-63-8 | 93   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN060824\  
Data File : BN031856.D  
Acq On : 08 Jun 2024 13:14  
Operator : MA/JU  
Sample : P2634-15  
Misc :  
ALS Vial : 7 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
BHBX9

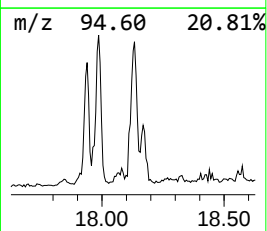
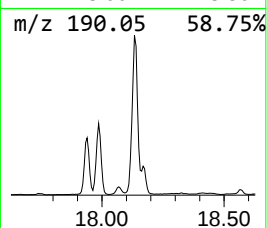
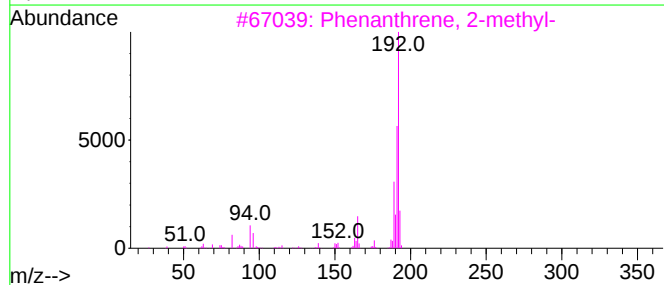
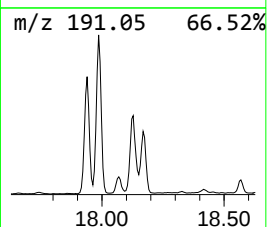
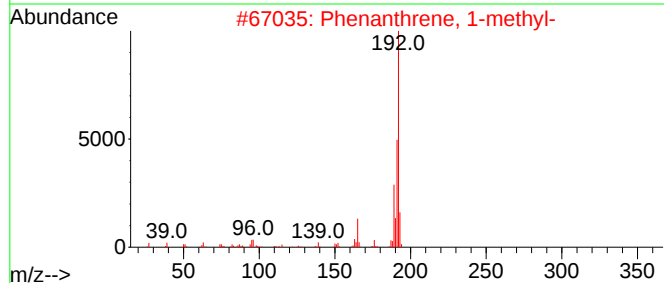
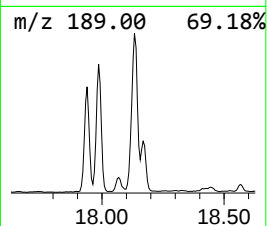
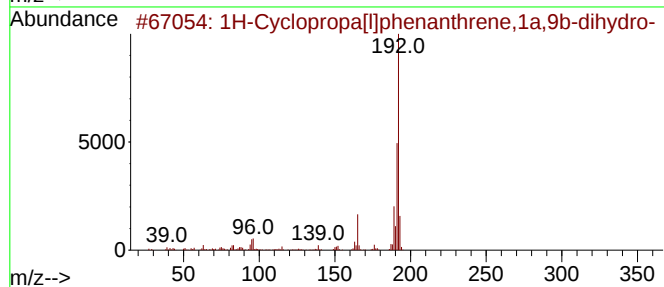
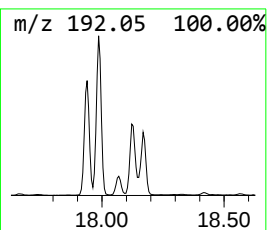
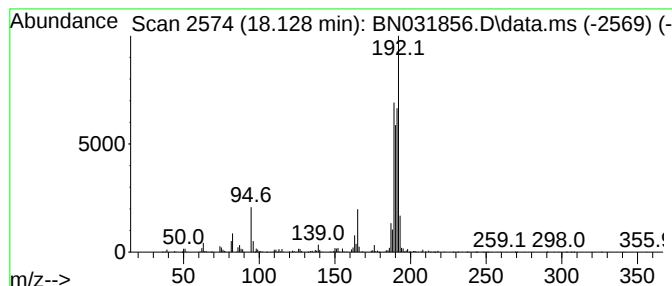
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Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 10 1H-Cyclopropa[l]phenanthren... Concentration Rank 4

| R.T.   | EstConc    | Area    | Relative to ISTD | R.T.   |
|--------|------------|---------|------------------|--------|
| 18.128 | 5.90 ng/ul | 1580260 | Phenanthrene-d10 | 17.063 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|---------|-------------|------|
| 1    |    |   | 1H-Cyclopropa[l]phenanthrene,1a,... | 192 | C15H12  | 000949-41-7 | 91   |
| 2    |    |   | Phenanthrene, 1-methyl-             | 192 | C15H12  | 000832-69-9 | 87   |
| 3    |    |   | Phenanthrene, 2-methyl-             | 192 | C15H12  | 002531-84-2 | 64   |
| 4    |    |   | Phenanthrene, 4-methyl-             | 192 | C15H12  | 000832-64-4 | 60   |
| 5    |    |   | Naphtho[2,3-b]norbornadiene         | 192 | C15H12  | 107426-38-0 | 60   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN060824\  
Data File : BN031856.D  
Acq On : 08 Jun 2024 13:14  
Operator : MA/JU  
Sample : P2634-15  
Misc :  
ALS Vial : 7 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
BHBX9

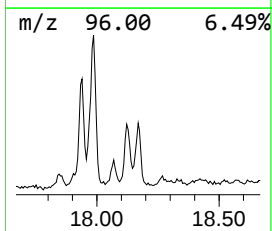
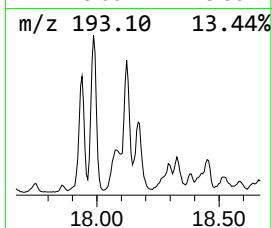
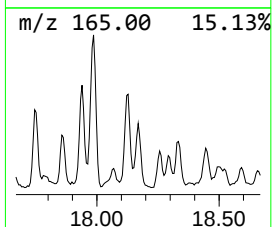
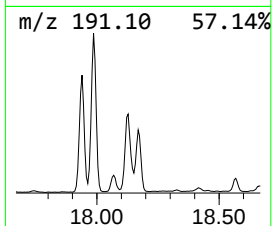
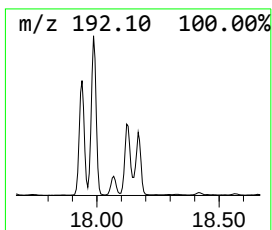
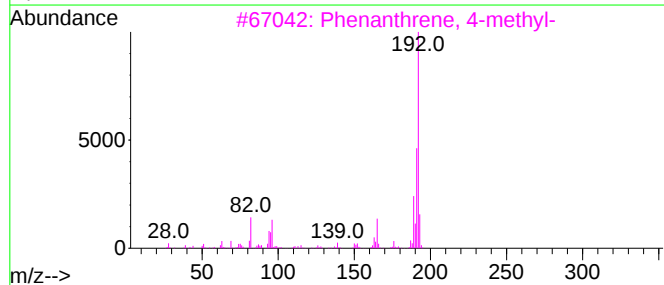
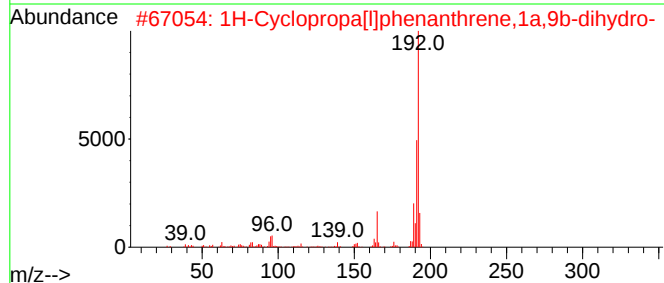
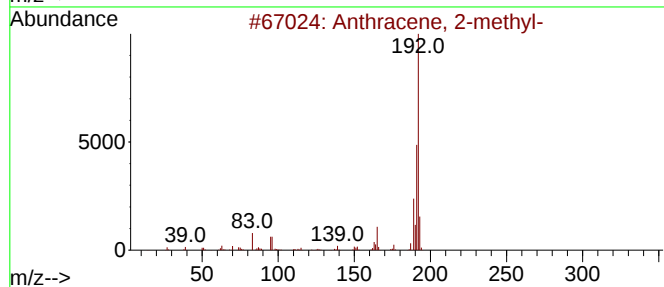
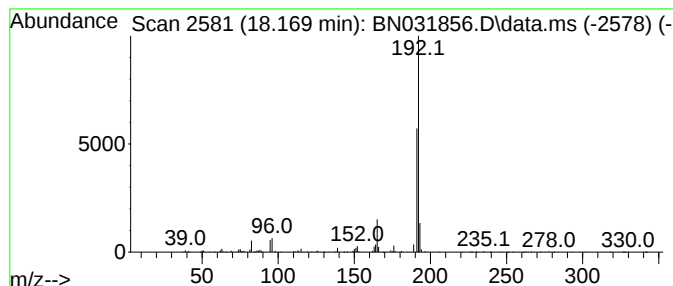
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Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 11 Anthracene, 2-methyl- Concentration Rank 17

| R.T.   | EstConc    | Area   | Relative to ISTD | R.T.   |
|--------|------------|--------|------------------|--------|
| 18.169 | 2.90 ng/ul | 775932 | Phenanthrene-d10 | 17.063 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|---------|-------------|------|
| 1    |    |   | Anthracene, 2-methyl-               | 192 | C15H12  | 000613-12-7 | 94   |
| 2    |    |   | 1H-Cyclopropa[1]phenanthrene,1a,... | 192 | C15H12  | 000949-41-7 | 94   |
| 3    |    |   | Phenanthrene, 4-methyl-             | 192 | C15H12  | 000832-64-4 | 91   |
| 4    |    |   | Anthracene, 9-methyl-               | 192 | C15H12  | 000779-02-2 | 91   |
| 5    |    |   | Anthracene, 9-methyl-               | 192 | C15H12  | 000779-02-2 | 91   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN060824\  
Data File : BN031856.D  
Acq On : 08 Jun 2024 13:14  
Operator : MA/JU  
Sample : P2634-15  
Misc :  
ALS Vial : 7 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
BHBX9

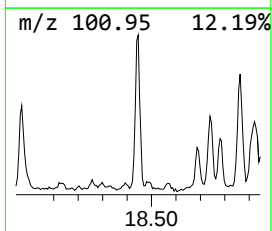
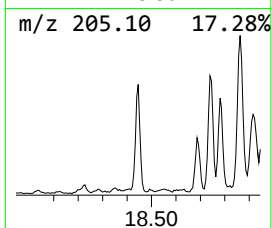
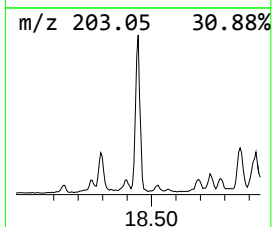
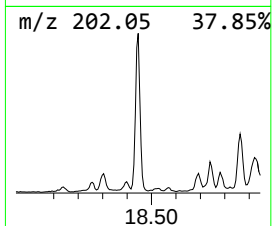
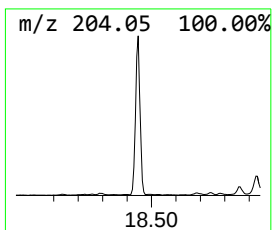
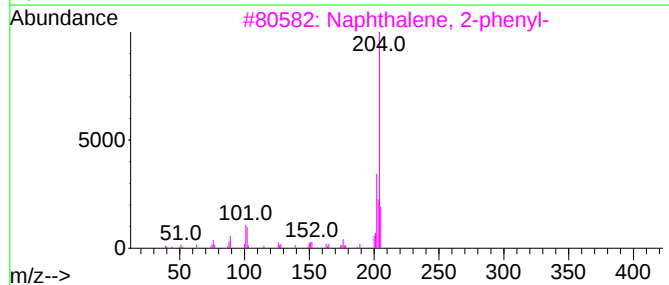
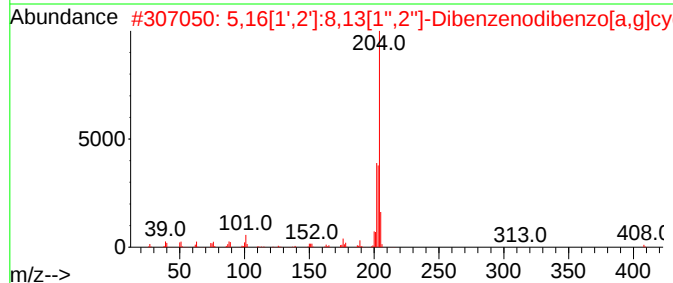
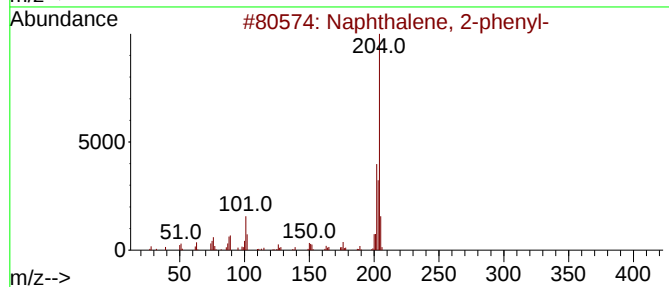
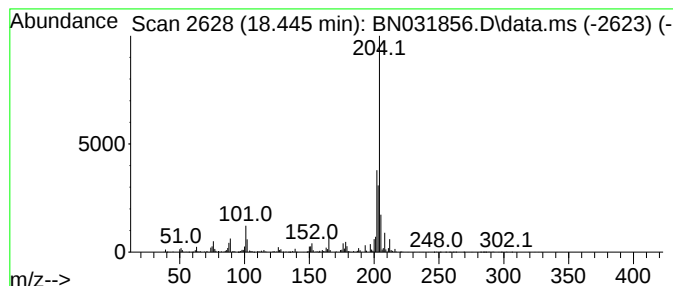
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Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 12 Naphthalene, 2-phenyl- Concentration Rank 10

| R.T.   | EstConc    | Area   | Relative to ISTD | R.T.   |
|--------|------------|--------|------------------|--------|
| 18.445 | 3.69 ng/ul | 988100 | Phenanthrene-d10 | 17.063 |

| Hit# | of 5 | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|------|-------------------------------------|-----|---------|-------------|------|
| 1    |      | Naphthalene, 2-phenyl-              | 204 | C16H12  | 000612-94-2 | 90   |
| 2    |      | 5,16[1',2']:8,13[1'',2'']-Dibenz... | 408 | C32H24  | 005672-97-9 | 87   |
| 3    |      | Naphthalene, 2-phenyl-              | 204 | C16H12  | 000612-94-2 | 86   |
| 4    |      | Naphthalene, 2-phenyl-              | 204 | C16H12  | 000612-94-2 | 81   |
| 5    |      | Tricyclo[8.2.2.2(4,7)]hexadeca-2... | 204 | C16H12  | 006572-60-7 | 76   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN060824\  
Data File : BN031856.D  
Acq On : 08 Jun 2024 13:14  
Operator : MA/JU  
Sample : P2634-15  
Misc :  
ALS Vial : 7 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
BHBX9

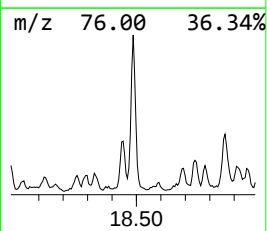
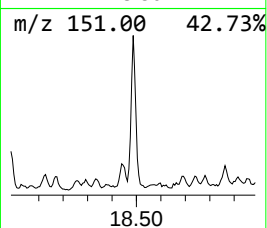
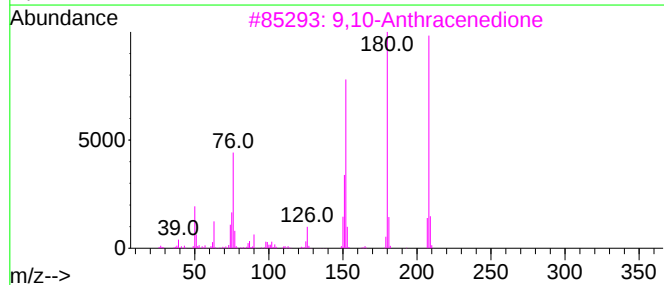
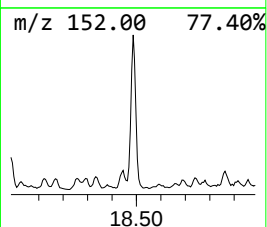
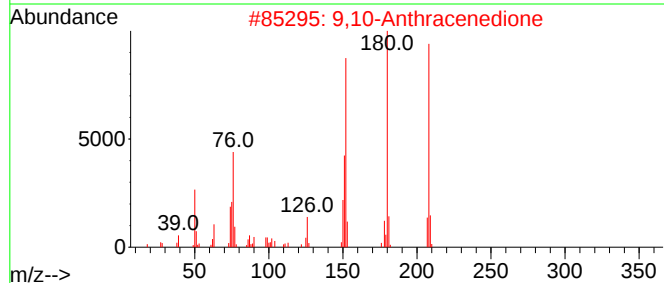
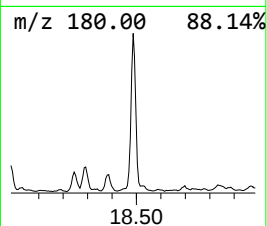
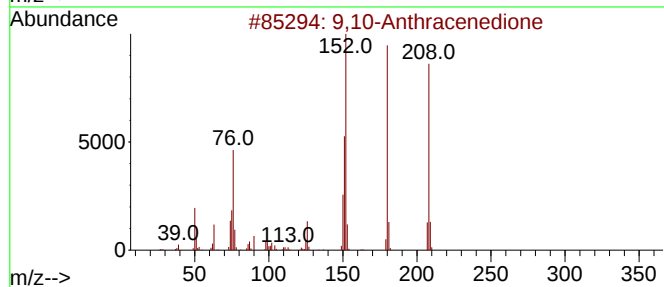
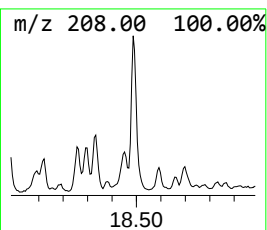
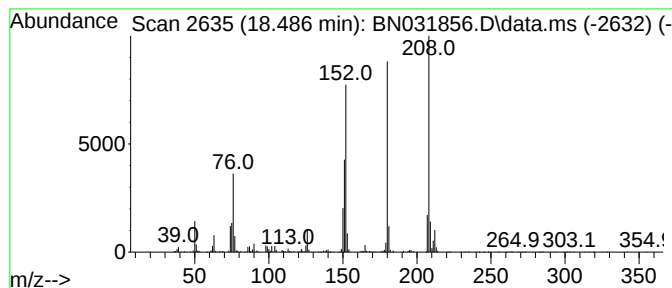
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Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 13 9,10-Anthracenedione Concentration Rank 19

| R.T.   | EstConc    | Area   | Relative to ISTD | R.T.   |
|--------|------------|--------|------------------|--------|
| 18.486 | 2.67 ng/ul | 715658 | Phenanthrene-d10 | 17.063 |

| Hit# | of                   | 5 | Tentative ID | MW  | MolForm | CAS#        | Qual |
|------|----------------------|---|--------------|-----|---------|-------------|------|
| 1    | 9,10-Anthracenedione |   |              | 208 | C14H8O2 | 000084-65-1 | 98   |
| 2    | 9,10-Anthracenedione |   |              | 208 | C14H8O2 | 000084-65-1 | 97   |
| 3    | 9,10-Anthracenedione |   |              | 208 | C14H8O2 | 000084-65-1 | 96   |
| 4    | 9H-Fluoren-9-one     |   |              | 180 | C13H8O  | 000486-25-9 | 91   |
| 5    | 9,10-Anthracenedione |   |              | 208 | C14H8O2 | 000084-65-1 | 89   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN060824\  
Data File : BN031856.D  
Acq On : 08 Jun 2024 13:14  
Operator : MA/JU  
Sample : P2634-15  
Misc :  
ALS Vial : 7 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
BHBX9

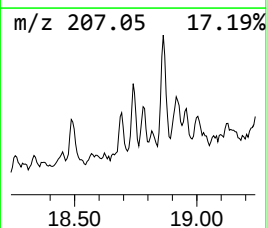
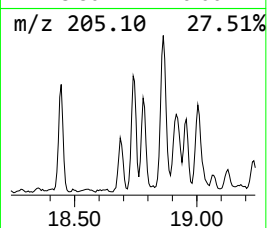
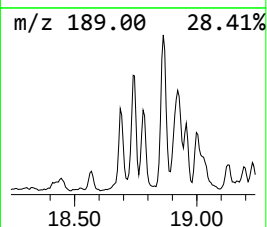
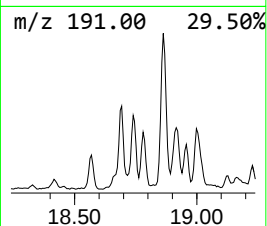
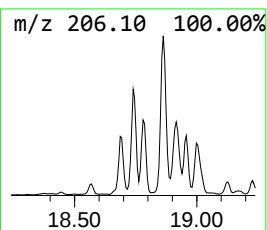
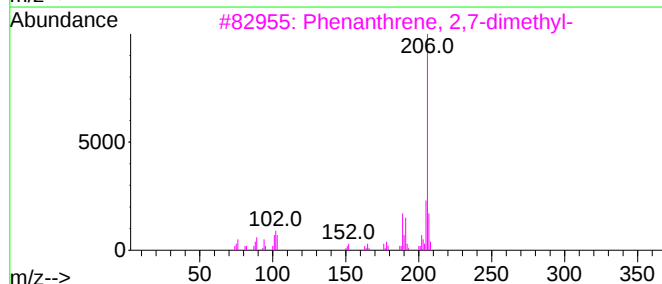
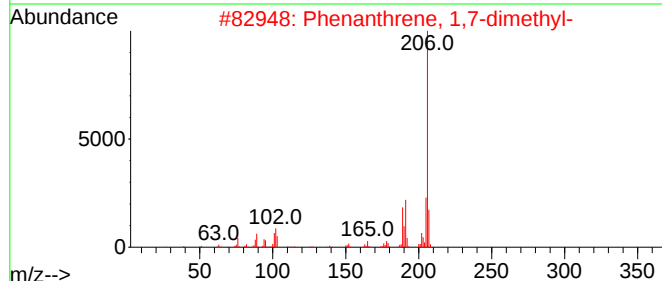
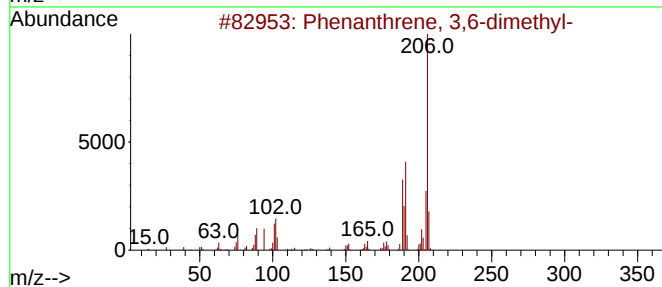
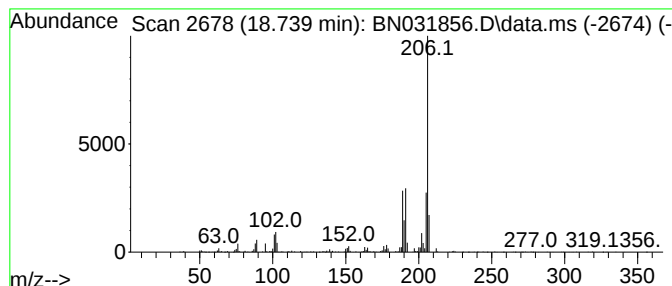
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Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 14 Phenanthrene, 3,6-dimethyl- Concentration Rank 24

| R.T.   | EstConc    | Area   | Relative to ISTD | R.T.   |
|--------|------------|--------|------------------|--------|
| 18.739 | 2.41 ng/ul | 646847 | Phenanthrene-d10 | 17.063 |

| Hit# | of | 5 | Tentative ID                | MW  | MolForm | CAS#        | Qual |
|------|----|---|-----------------------------|-----|---------|-------------|------|
| 1    |    |   | Phenanthrene, 3,6-dimethyl- | 206 | C16H14  | 001576-67-6 | 97   |
| 2    |    |   | Phenanthrene, 1,7-dimethyl- | 206 | C16H14  | 000483-87-4 | 93   |
| 3    |    |   | Phenanthrene, 2,7-dimethyl- | 206 | C16H14  | 001576-69-8 | 93   |
| 4    |    |   | di-p-Tolylacetylene         | 206 | C16H14  | 002789-88-0 | 91   |
| 5    |    |   | Phenanthrene, 2,7-dimethyl- | 206 | C16H14  | 001576-69-8 | 90   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN060824\  
Data File : BN031856.D  
Acq On : 08 Jun 2024 13:14  
Operator : MA/JU  
Sample : P2634-15  
Misc :  
ALS Vial : 7 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
BHBX9

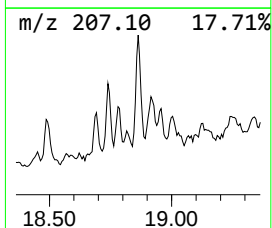
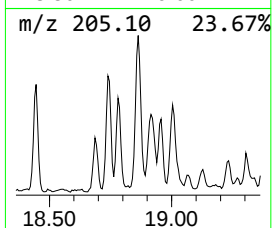
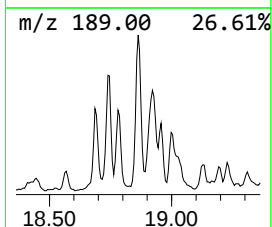
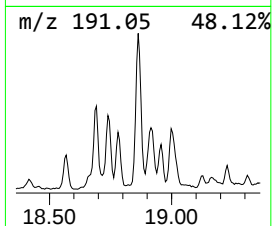
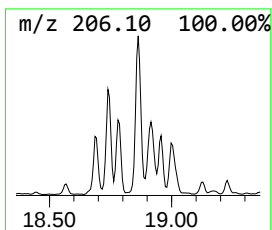
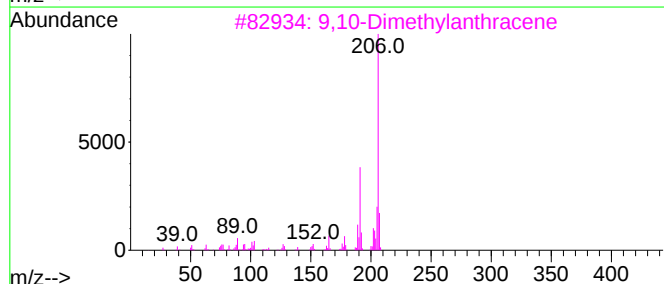
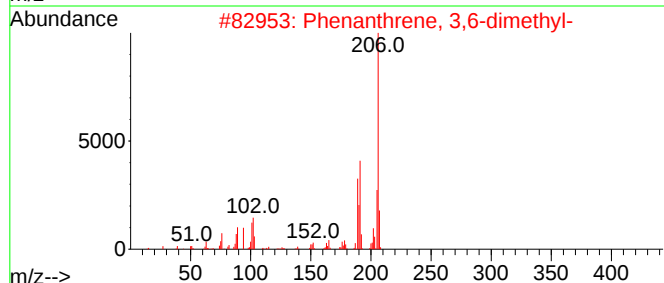
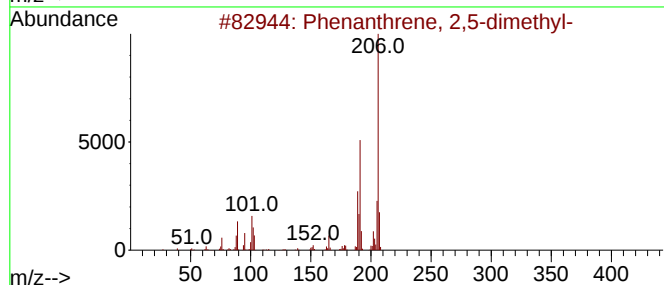
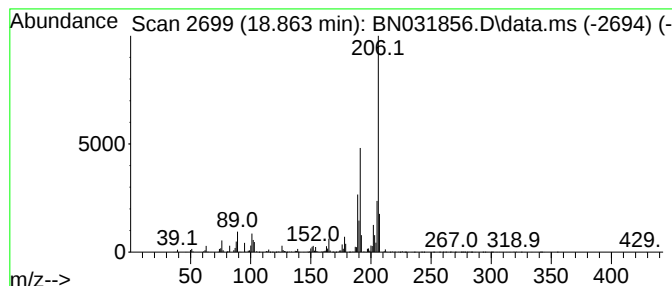
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Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 15 Phenanthrene, 2,5-dimethyl- Concentration Rank 8

| R.T.   | EstConc    | Area    | Relative to ISTD | R.T.   |
|--------|------------|---------|------------------|--------|
| 18.863 | 4.45 ng/ul | 1192940 | Phenanthrene-d10 | 17.063 |

| Hit# | of | 5 | Tentative ID                | MW  | MolForm | CAS#        | Qual |
|------|----|---|-----------------------------|-----|---------|-------------|------|
| 1    |    |   | Phenanthrene, 2,5-dimethyl- | 206 | C16H14  | 003674-66-6 | 98   |
| 2    |    |   | Phenanthrene, 3,6-dimethyl- | 206 | C16H14  | 001576-67-6 | 93   |
| 3    |    |   | 9,10-Dimethylanthracene     | 206 | C16H14  | 000781-43-1 | 93   |
| 4    |    |   | 9,10-Dimethylanthracene     | 206 | C16H14  | 000781-43-1 | 93   |
| 5    |    |   | Phenanthrene, 2,3-dimethyl- | 206 | C16H14  | 003674-65-5 | 93   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN060824\  
Data File : BN031856.D  
Acq On : 08 Jun 2024 13:14  
Operator : MA/JU  
Sample : P2634-15  
Misc :  
ALS Vial : 7 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
BHBX9

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_N\Methods\SFAM-EPA-BN052924.MA.M  
Quant Title : SVOA CALIBRATION

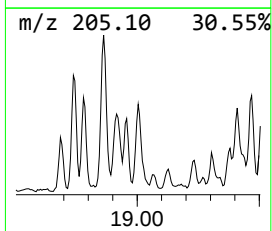
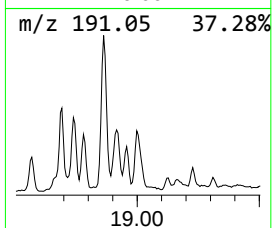
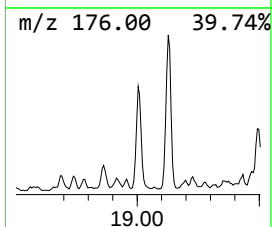
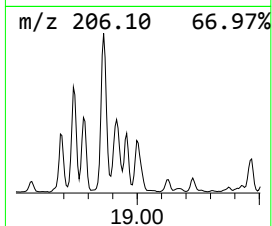
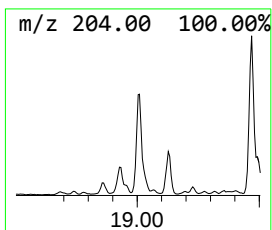
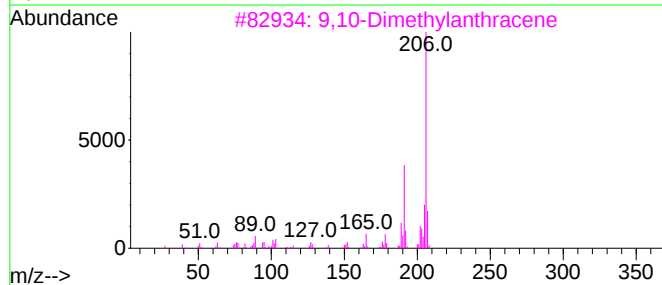
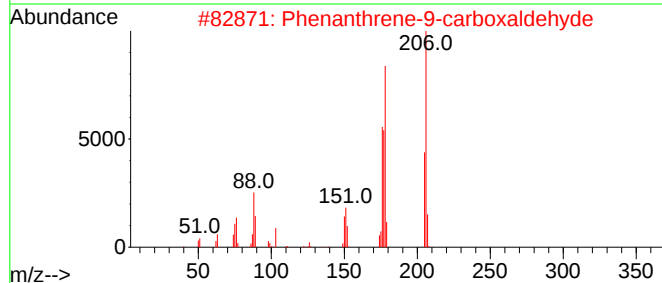
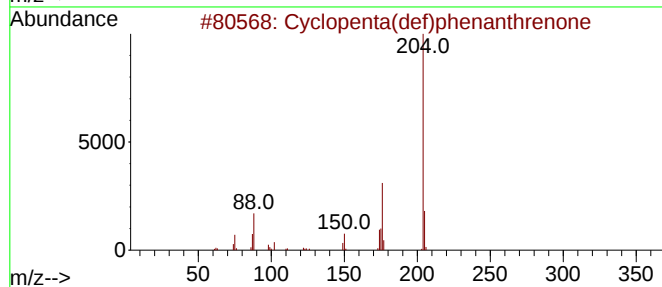
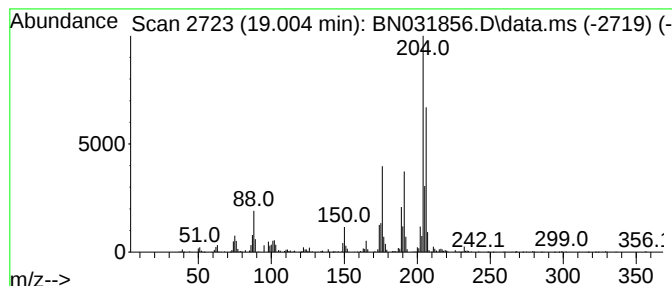
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TIC Integration Parameters: LSCINT.P

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Peak Number 16 Cyclopenta(def)phenanthrene Concentration Rank 15

| R.T.   | EstConc    | Area   | Relative to ISTD | R.T.   |
|--------|------------|--------|------------------|--------|
| 19.004 | 3.21 ng/ul | 860082 | Phenanthrene-d10 | 17.063 |

| Hit# | of | 5 | Tentative ID                  | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------------|-----|---------|-------------|------|
| 1    |    |   | Cyclopenta(def)phenanthrene   | 204 | C15H8O  | 005737-13-3 | 95   |
| 2    |    |   | Phenanthrene-9-carboxaldehyde | 206 | C15H10O | 004707-71-5 | 43   |
| 3    |    |   | 9,10-Dimethylantracene        | 206 | C16H14  | 000781-43-1 | 38   |
| 4    |    |   | 9,10-Dimethylantracene        | 206 | C16H14  | 000781-43-1 | 38   |
| 5    |    |   | 9,10-Dimethylantracene        | 206 | C16H14  | 000781-43-1 | 38   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN060824\  
Data File : BN031856.D  
Acq On : 08 Jun 2024 13:14  
Operator : MA/JU  
Sample : P2634-15  
Misc :  
ALS Vial : 7 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
BHBX9

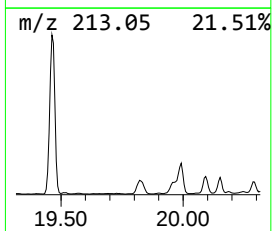
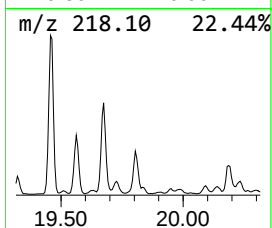
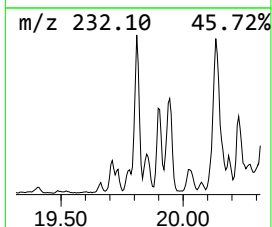
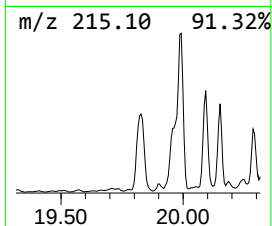
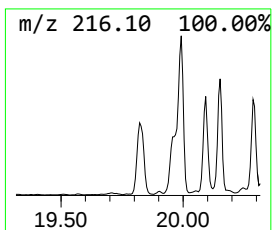
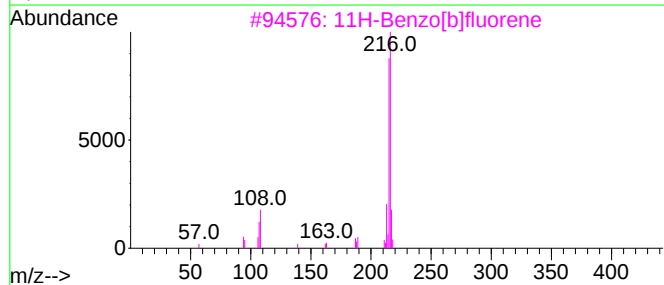
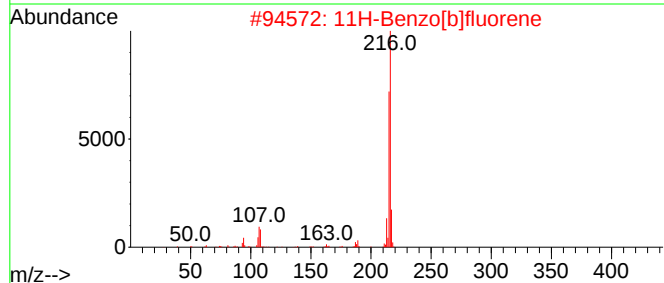
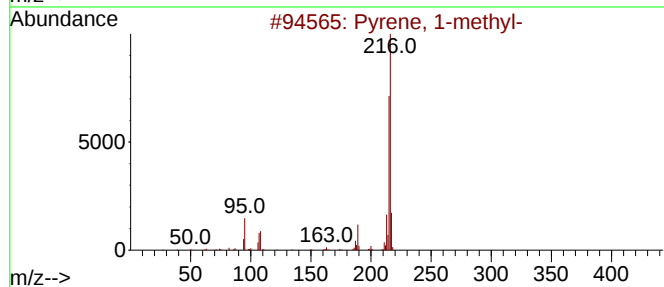
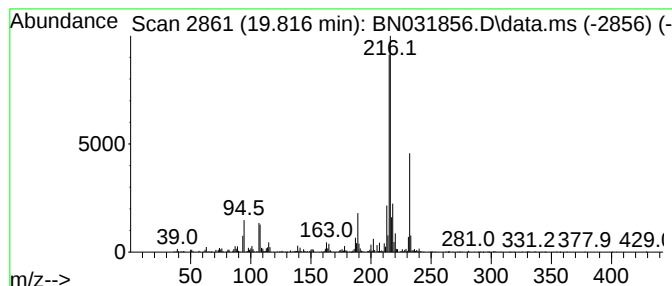
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Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 17 Pyrene, 1-methyl- Concentration Rank 13

| R.T.   | EstConc    | Area    | Relative to ISTD | R.T.   |
|--------|------------|---------|------------------|--------|
| 19.816 | 3.24 ng/ul | 1264840 | Chrysene-d12     | 21.304 |

| Hit# | of | 5 | Tentative ID         | MW  | MolForm | CAS#        | Qual |
|------|----|---|----------------------|-----|---------|-------------|------|
| 1    |    |   | Pyrene, 1-methyl-    | 216 | C17H12  | 002381-21-7 | 96   |
| 2    |    |   | 11H-Benzo[b]fluorene | 216 | C17H12  | 000243-17-4 | 94   |
| 3    |    |   | 11H-Benzo[b]fluorene | 216 | C17H12  | 000243-17-4 | 92   |
| 4    |    |   | Pyrene, 1-methyl-    | 216 | C17H12  | 002381-21-7 | 90   |
| 5    |    |   | Pyrene, 1-methyl-    | 216 | C17H12  | 002381-21-7 | 86   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN060824\  
Data File : BN031856.D  
Acq On : 08 Jun 2024 13:14  
Operator : MA/JU  
Sample : P2634-15  
Misc :  
ALS Vial : 7 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
BHBX9

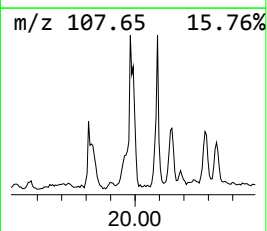
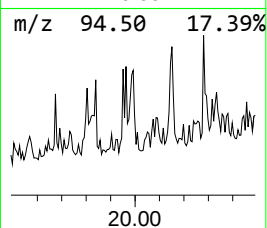
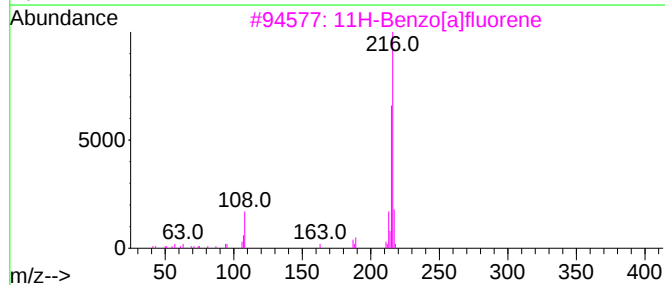
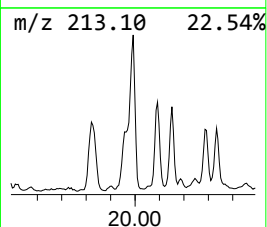
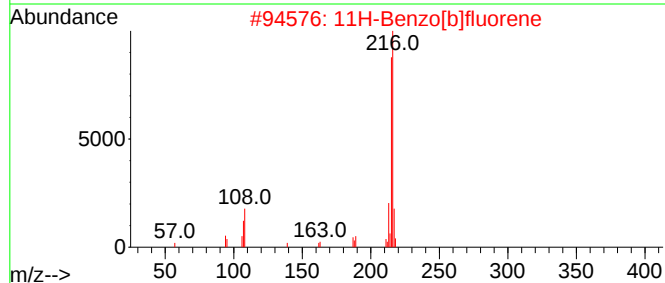
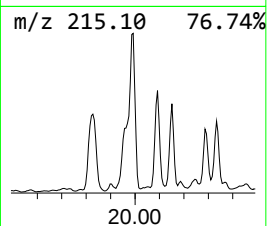
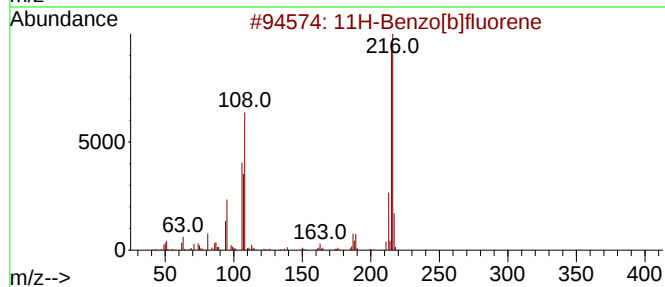
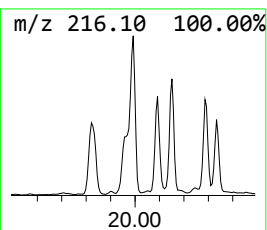
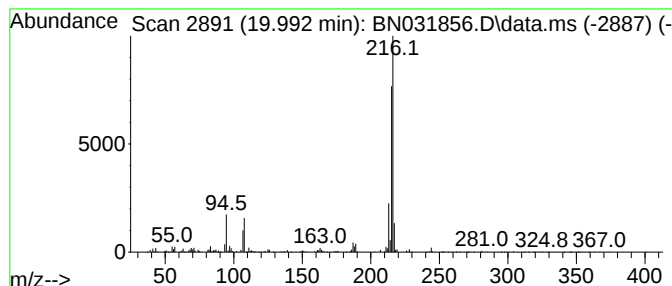
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Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 18 11H-Benzo[b]fluorene Concentration Rank 26

| R.T.   | EstConc    | Area   | Relative to ISTD | R.T.   |
|--------|------------|--------|------------------|--------|
| 19.992 | 2.18 ng/ul | 849534 | Chrysene-d12     | 21.304 |

| Hit# | of | 5 | Tentative ID         | MW  | MolForm | CAS#        | Qual |
|------|----|---|----------------------|-----|---------|-------------|------|
| 1    |    |   | 11H-Benzo[b]fluorene | 216 | C17H12  | 000243-17-4 | 91   |
| 2    |    |   | 11H-Benzo[b]fluorene | 216 | C17H12  | 000243-17-4 | 91   |
| 3    |    |   | 11H-Benzo[a]fluorene | 216 | C17H12  | 000238-84-6 | 90   |
| 4    |    |   | 11H-Benzo[b]fluorene | 216 | C17H12  | 000243-17-4 | 87   |
| 5    |    |   | 11H-Benzo[a]fluorene | 216 | C17H12  | 000238-84-6 | 83   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN060824\  
Data File : BN031856.D  
Acq On : 08 Jun 2024 13:14  
Operator : MA/JU  
Sample : P2634-15  
Misc :  
ALS Vial : 7 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
BHBX9

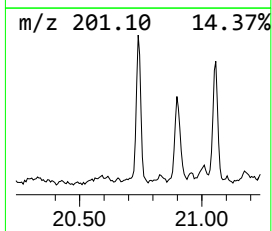
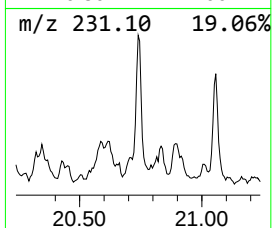
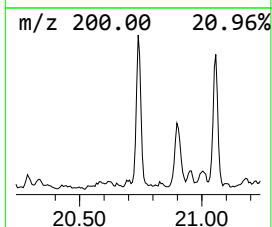
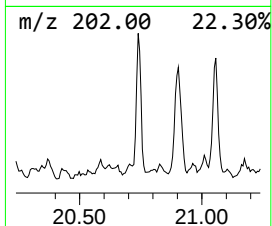
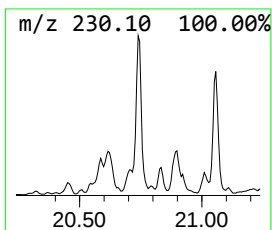
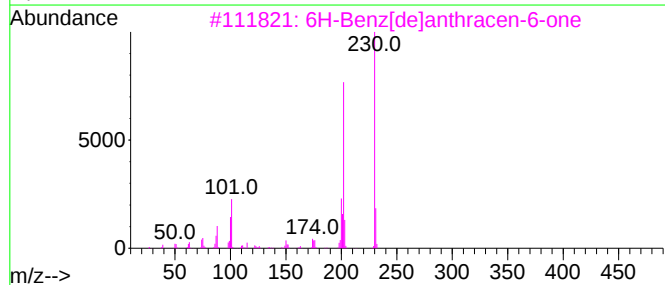
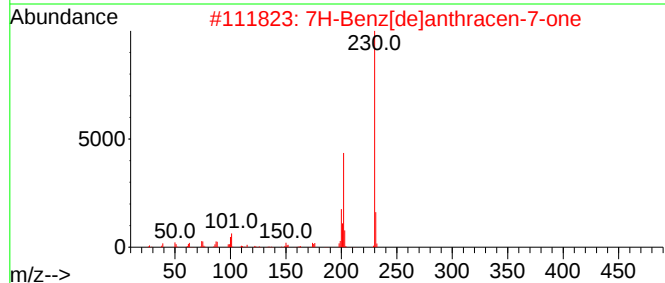
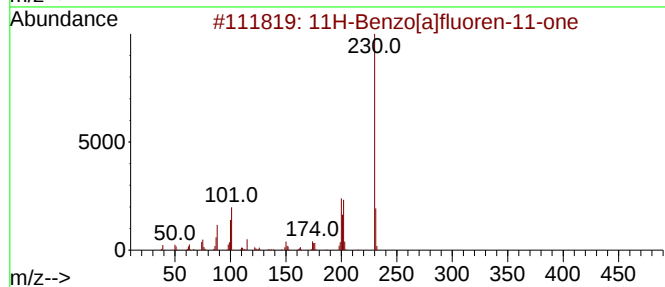
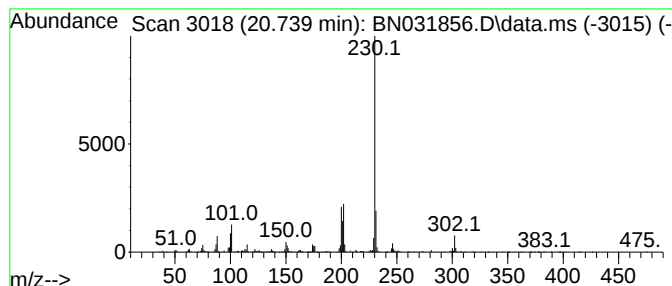
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Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 19 11H-Benzo[a]fluoren-11-one Concentration Rank 23

| R.T.   | EstConc    | Area   | Relative to ISTD | R.T.   |
|--------|------------|--------|------------------|--------|
| 20.739 | 2.42 ng/ul | 945004 | Chrysene-d12     | 21.304 |

| Hit# | of | 5 | Tentative ID               | MW  | MolForm | CAS#        | Qual |
|------|----|---|----------------------------|-----|---------|-------------|------|
| 1    |    |   | 11H-Benzo[a]fluoren-11-one | 230 | C17H10O | 000479-79-8 | 98   |
| 2    |    |   | 7H-Benz[de]anthracen-7-one | 230 | C17H10O | 000082-05-3 | 90   |
| 3    |    |   | 6H-Benz[de]anthracen-6-one | 230 | C17H10O | 080252-14-8 | 90   |
| 4    |    |   | 7H-Benz[de]anthracen-7-one | 230 | C17H10O | 000082-05-3 | 87   |
| 5    |    |   | 7H-Benz[de]anthracen-7-one | 230 | C17H10O | 000082-05-3 | 87   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN060824\  
Data File : BN031856.D  
Acq On : 08 Jun 2024 13:14  
Operator : MA/JU  
Sample : P2634-15  
Misc :  
ALS Vial : 7 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
BHBX9

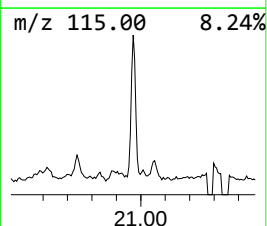
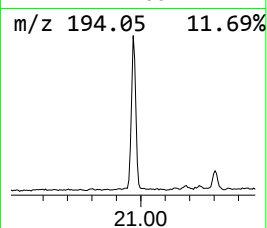
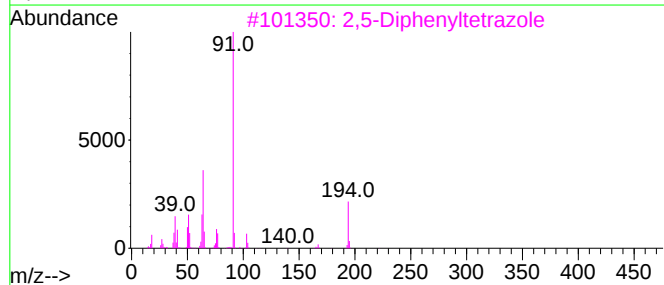
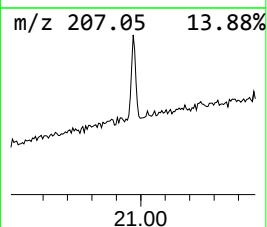
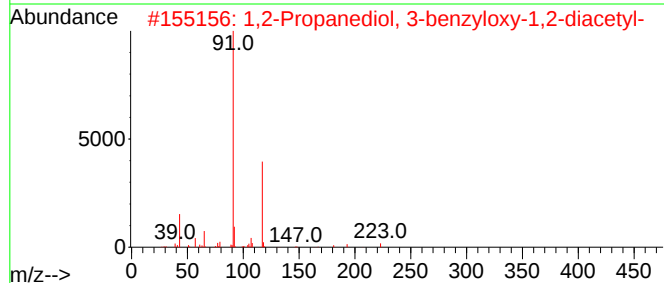
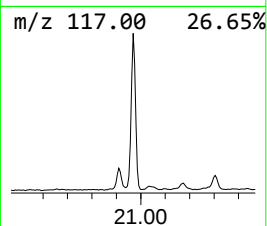
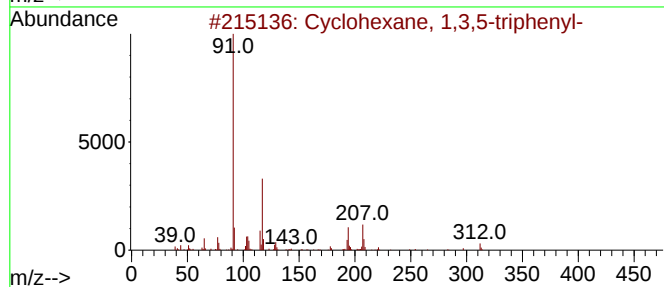
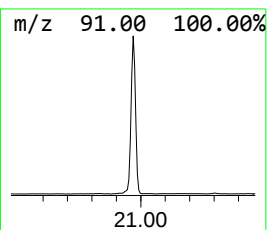
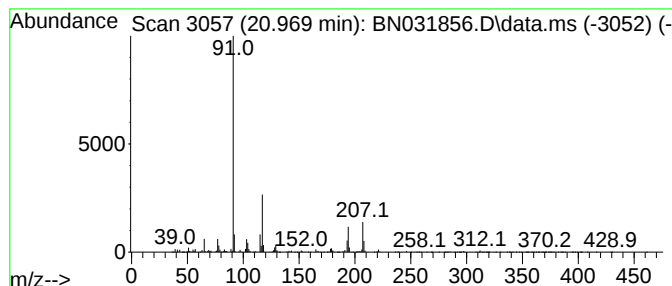
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Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 20 Cyclohexane, 1,3,5-triphenyl- Concentration Rank 3

| R.T.   | EstConc    | Area    | Relative to ISTD | R.T.   |
|--------|------------|---------|------------------|--------|
| 20.968 | 6.96 ng/ul | 2714290 | Chrysene-d12     | 21.304 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm   | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|-----------|--------------|------|
| 1    |    |   | Cyclohexane, 1,3,5-triphenyl-       | 312 | C24H24    | 028336-57-4  | 83   |
| 2    |    |   | 1,2-Propanediol, 3-benzyloxy-1,2... | 266 | C14H18O5  | 013754-10-4  | 32   |
| 3    |    |   | 2,5-Diphenyltetrazole               | 222 | C13H10N4  | 018039-45-7  | 23   |
| 4    |    |   | (R)-(1-Cyclopentylpropane-1,3-di... | 264 | C20H24    | 1402851-74-4 | 16   |
| 5    |    |   | Carbonic acid, monoamide, N-benz... | 207 | C12H17NO2 | 1000451-48-5 | 16   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN060824\  
Data File : BN031856.D  
Acq On : 08 Jun 2024 13:14  
Operator : MA/JU  
Sample : P2634-15  
Misc :  
ALS Vial : 7 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
BHBX9

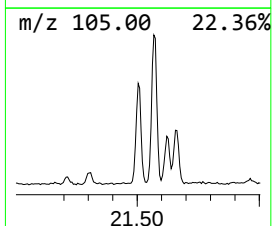
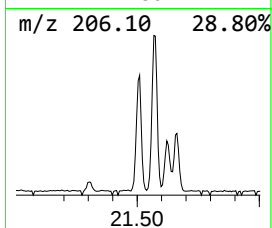
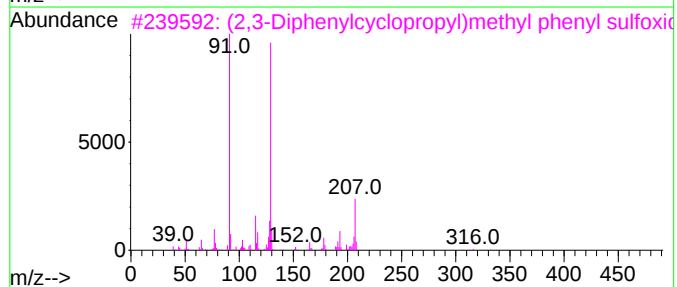
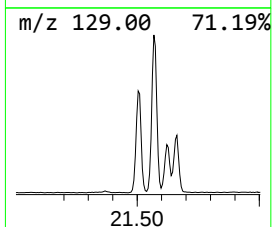
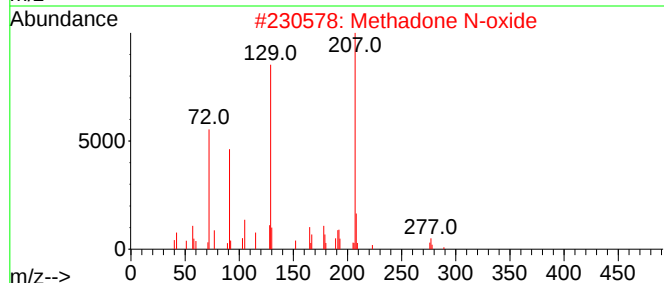
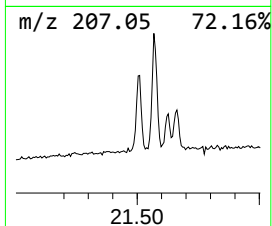
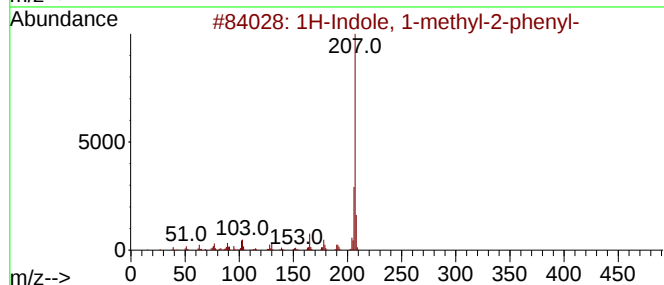
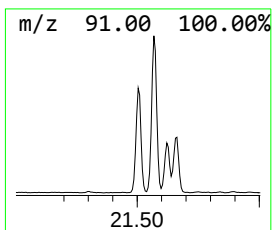
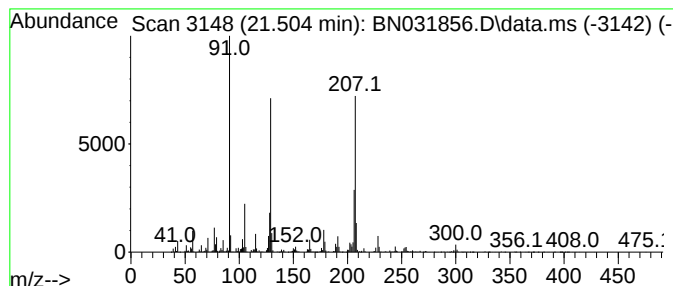
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Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 21 1H-Indole, 1-methyl-2-phenyl- Concentration Rank 2

| R.T.   | EstConc    | Area    | Relative to ISTD | R.T.   |
|--------|------------|---------|------------------|--------|
| 21.504 | 7.02 ng/ul | 2737540 | Chrysene-d12     | 21.304 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm   | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|-----------|--------------|------|
| 1    |    |   | 1H-Indole, 1-methyl-2-phenyl-       | 207 | C15H13N   | 003558-24-5  | 55   |
| 2    |    |   | Methadone N-oxide                   | 325 | C21H27NO2 | 1000120-80-7 | 47   |
| 3    |    |   | (2,3-Diphenylcyclopropyl)methyl ... | 332 | C22H20OS  | 131758-71-9  | 47   |
| 4    |    |   | 1H-Indole, 2-methyl-3-phenyl-       | 207 | C15H13N   | 004757-69-1  | 46   |
| 5    |    |   | 1H-Indole, 2-methyl-3-phenyl-       | 207 | C15H13N   | 004757-69-1  | 41   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN060824\  
Data File : BN031856.D  
Acq On : 08 Jun 2024 13:14  
Operator : MA/JU  
Sample : P2634-15  
Misc :  
ALS Vial : 7 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
BHBX9

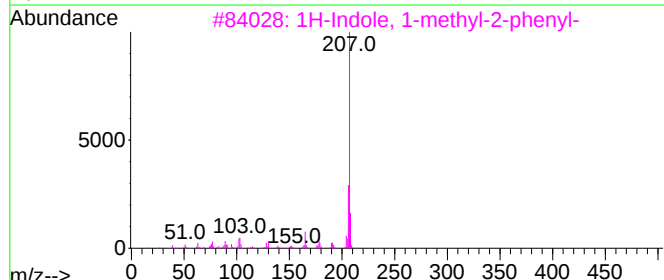
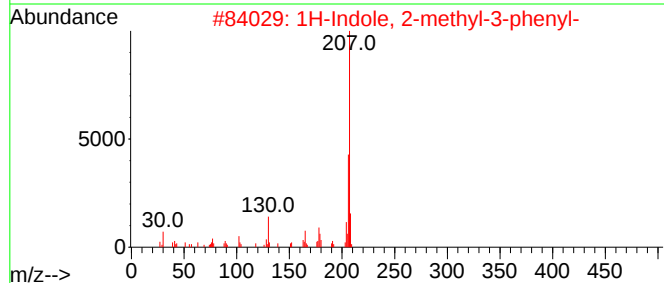
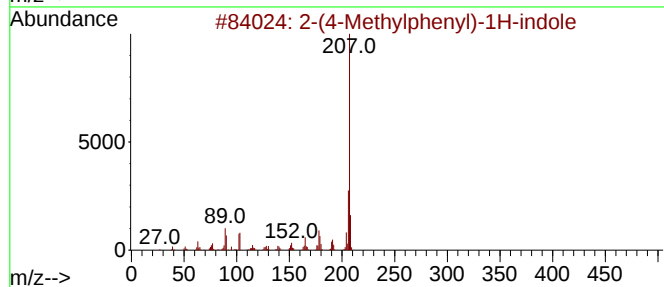
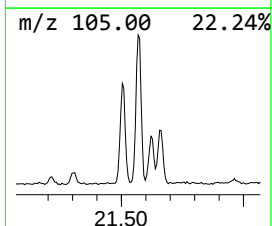
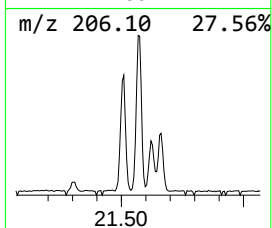
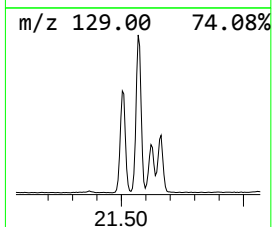
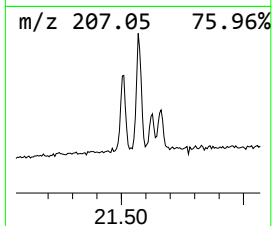
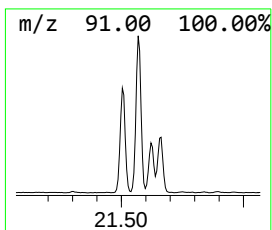
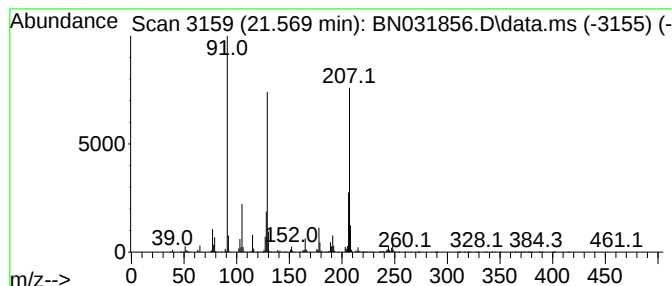
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Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 22 unknown-02 Concentration Rank 1

| R.T.   | EstConc    | Area    | Relative to ISTD | R.T.   |
|--------|------------|---------|------------------|--------|
| 21.569 | 8.17 ng/ul | 3183340 | Chrysene-d12     | 21.304 |

| Hit# | of                            | 5 | Tentative ID | MW  | MolForm | CAS#         | Qual |
|------|-------------------------------|---|--------------|-----|---------|--------------|------|
| 1    | 2-(4-Methylphenyl)-1H-indole  |   |              | 207 | C15H13N | 055577-25-8  | 46   |
| 2    | 1H-Indole, 2-methyl-3-phenyl- |   |              | 207 | C15H13N | 004757-69-1  | 43   |
| 3    | 1H-Indole, 1-methyl-2-phenyl- |   |              | 207 | C15H13N | 003558-24-5  | 42   |
| 4    | 5-Methyl-2-phenylindolizine   |   |              | 207 | C15H13N | 036944-99-7  | 38   |
| 5    | 1-benzylindole                |   |              | 207 | C15H13N | 1000334-31-3 | 38   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN060824\  
Data File : BN031856.D  
Acq On : 08 Jun 2024 13:14  
Operator : MA/JU  
Sample : P2634-15  
Misc :  
ALS Vial : 7 Sample Multiplier: 1

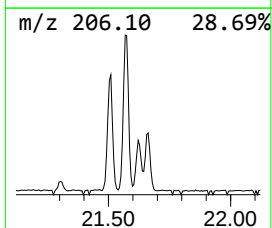
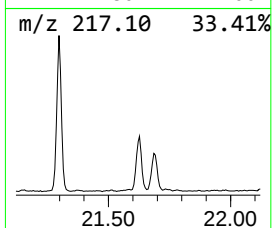
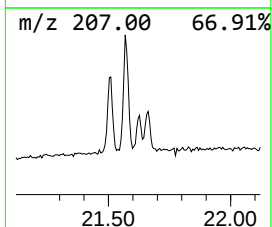
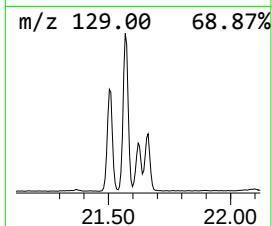
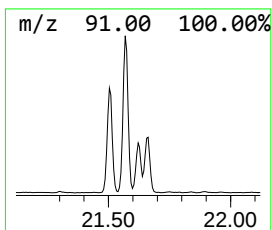
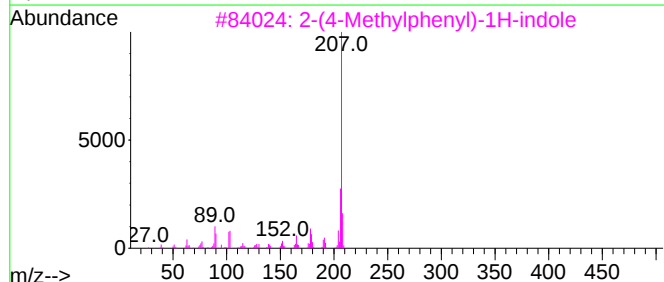
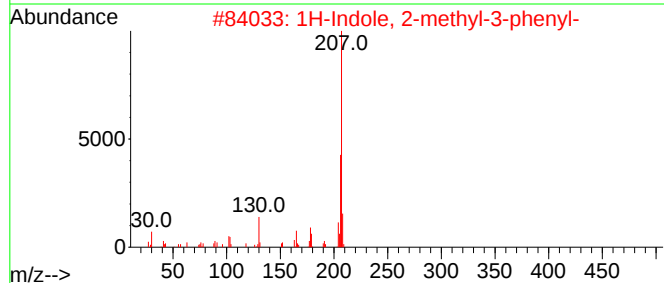
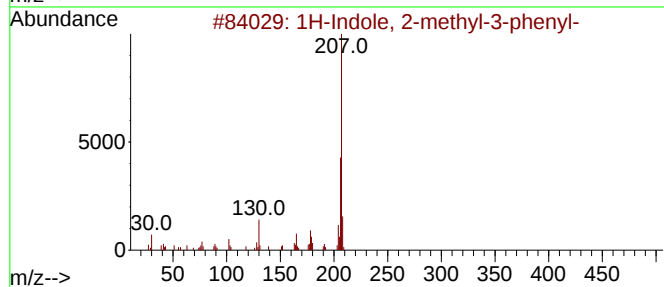
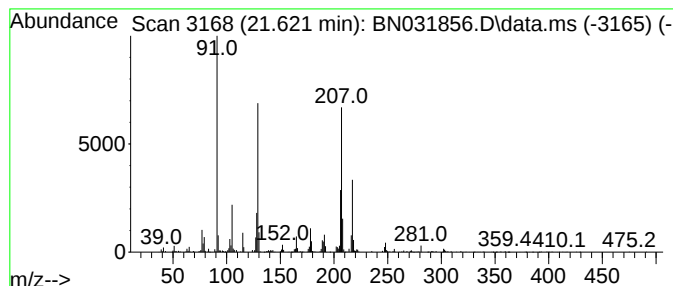
Instrument :  
BNA\_N  
ClientSampleId :  
BHBX9

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_N\Methods\SFAM-EPA-BN052924.MA.M  
Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 23 unknown-03 Concentration Rank 20

| R.T.   | EstConc    | Area   | Relative to ISTD              |     |         | R.T.        |      |
|--------|------------|--------|-------------------------------|-----|---------|-------------|------|
| 21.621 | 2.50 ng/ul | 972573 | Chrysene-d12                  |     |         | 21.304      |      |
| Hit#   | of         | 5      | Tentative ID                  | MW  | MolForm | CAS#        | Qual |
| 1      |            |        | 1H-Indole, 2-methyl-3-phenyl- | 207 | C15H13N | 004757-69-1 | 46   |
| 2      |            |        | 1H-Indole, 2-methyl-3-phenyl- | 207 | C15H13N | 004757-69-1 | 45   |
| 3      |            |        | 2-(4-Methylphenyl)-1H-indole  | 207 | C15H13N | 055577-25-8 | 42   |
| 4      |            |        | 1H-Indole, 1-methyl-2-phenyl- | 207 | C15H13N | 003558-24-5 | 38   |
| 5      |            |        | 1H-Indole, 5-methyl-2-phenyl- | 207 | C15H13N | 013228-36-9 | 38   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN060824\  
Data File : BN031856.D  
Acq On : 08 Jun 2024 13:14  
Operator : MA/JU  
Sample : P2634-15  
Misc :  
ALS Vial : 7 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
BHBX9

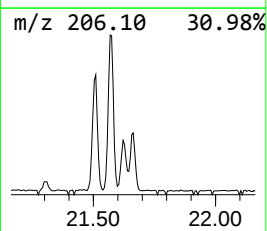
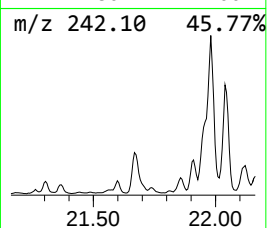
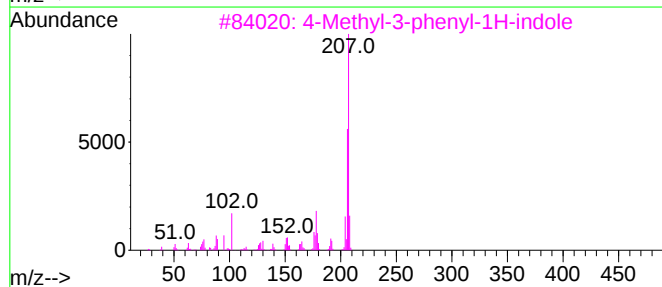
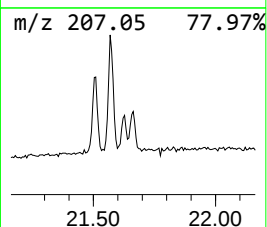
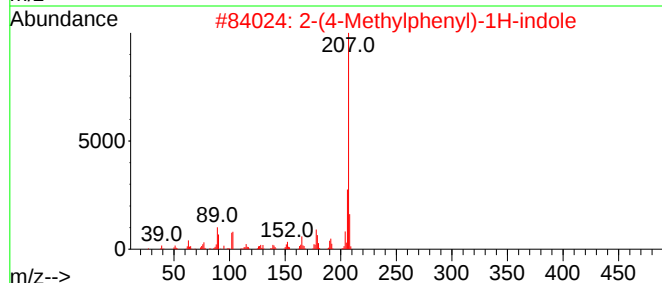
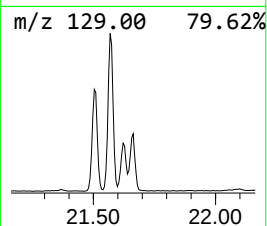
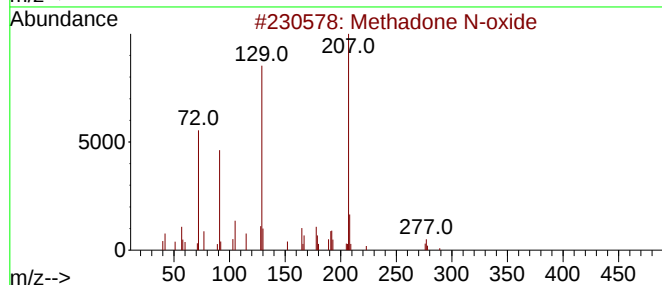
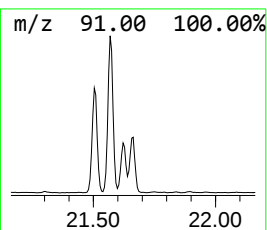
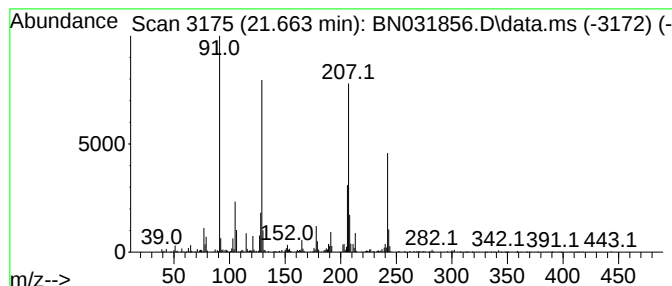
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Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 24 unknown-04 Concentration Rank 14

| R.T.   | EstConc    | Area    | Relative to ISTD | R.T.   |
|--------|------------|---------|------------------|--------|
| 21.663 | 3.21 ng/ul | 1252400 | Chrysene-d12     | 21.304 |

| Hit# | of | 5 | Tentative ID                 | MW  | MolForm    | CAS#         | Qual |
|------|----|---|------------------------------|-----|------------|--------------|------|
| 1    |    |   | Methadone N-oxide            | 325 | C21H27NO2  | 1000120-80-7 | 47   |
| 2    |    |   | 2-(4-Methylphenyl)-1H-indole | 207 | C15H13N    | 055577-25-8  | 46   |
| 3    |    |   | 4-Methyl-3-phenyl-1H-indole  | 207 | C15H13N    | 180271-54-9  | 38   |
| 4    |    |   | 5-Methyl-2-phenylindolizine  | 207 | C15H13N    | 036944-99-7  | 38   |
| 5    |    |   | Medazepam                    | 270 | C16H15ClN2 | 002898-12-6  | 38   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN060824\  
Data File : BN031856.D  
Acq On : 08 Jun 2024 13:14  
Operator : MA/JU  
Sample : P2634-15  
Misc :  
ALS Vial : 7 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
BHBX9

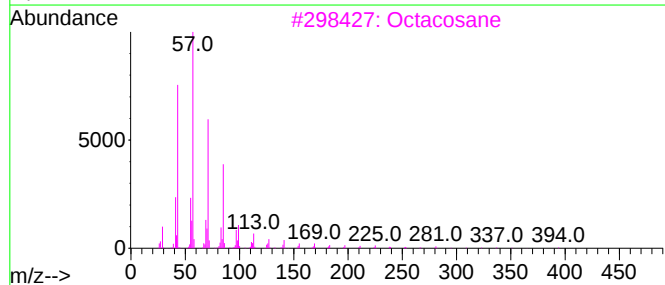
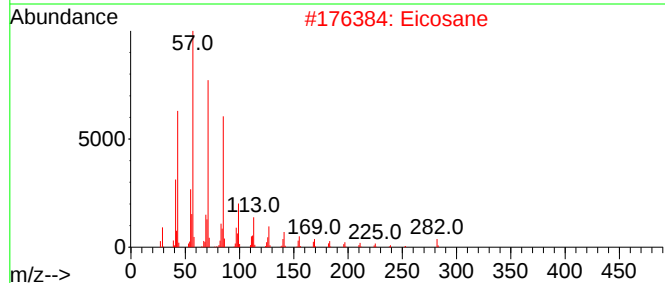
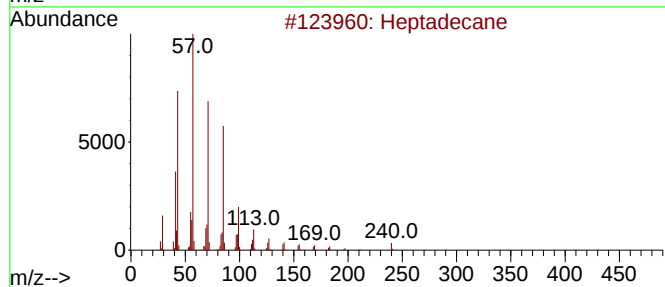
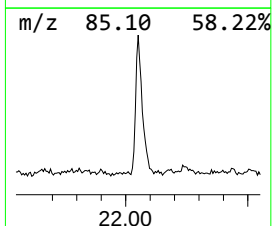
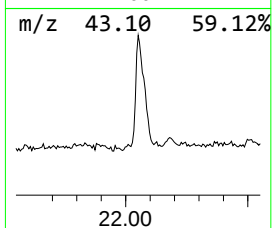
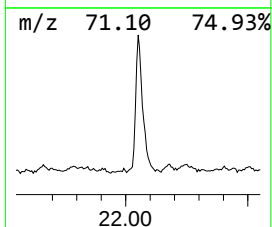
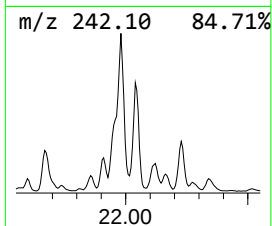
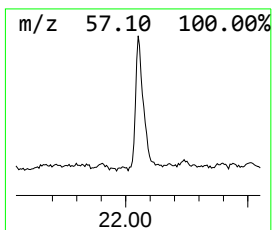
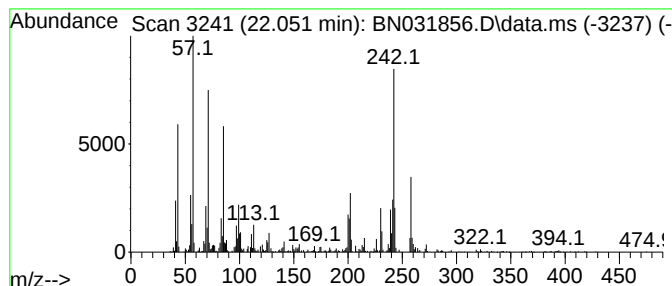
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Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 25 (DEL) Alkane: Straight-Chai... Concentration Rank 16

| R.T.      | EstConc             | Area    | Relative to ISTD |         | R.T.        |      |
|-----------|---------------------|---------|------------------|---------|-------------|------|
| 22.051    | 2.99 ng/ul          | 1166930 | Chrysene-d12     |         | 21.304      |      |
| Hit# of 5 | Tentative ID        |         | MW               | MolForm | CAS#        | Qual |
| 1         | Heptadecane         |         | 240              | C17H36  | 000629-78-7 | 80   |
| 2         | Eicosane            |         | 282              | C20H42  | 000112-95-8 | 80   |
| 3         | Octacosane          |         | 394              | C28H58  | 000630-02-4 | 72   |
| 4         | Hexadecane          |         | 226              | C16H34  | 000544-76-3 | 59   |
| 5         | Chrysene, 1-methyl- |         | 242              | C19H14  | 003351-28-8 | 55   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN060824\  
Data File : BN031856.D  
Acq On : 08 Jun 2024 13:14  
Operator : MA/JU  
Sample : P2634-15  
Misc :  
ALS Vial : 7 Sample Multiplier: 1

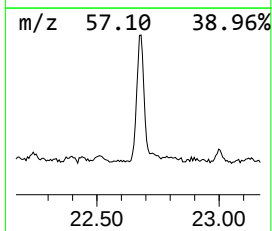
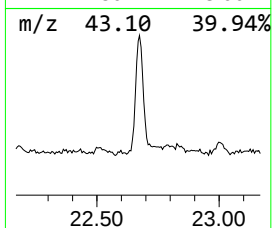
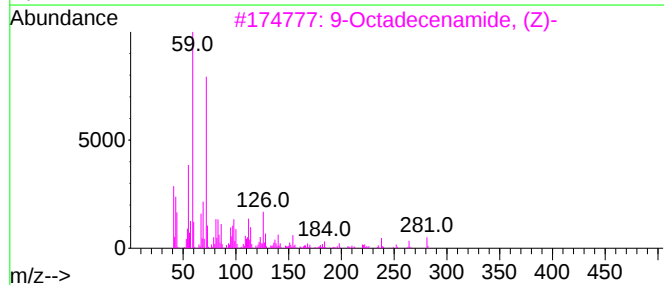
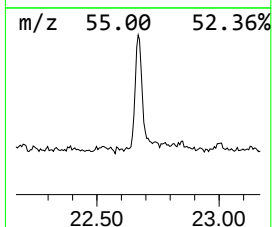
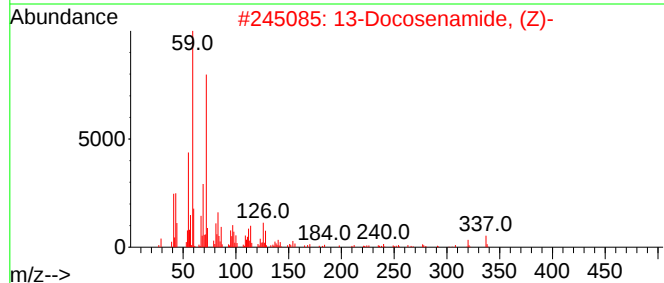
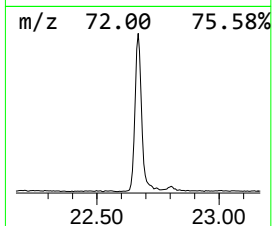
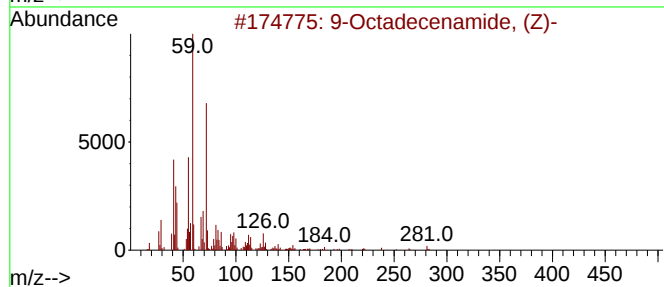
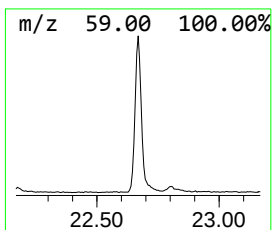
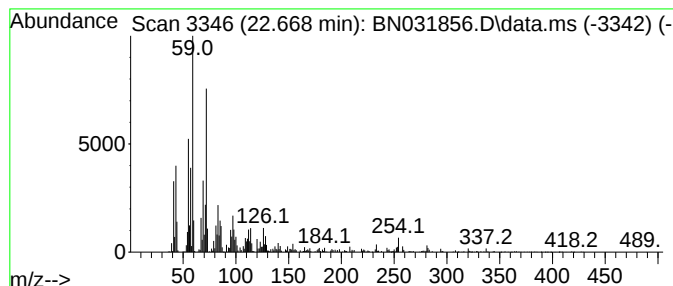
Instrument :  
BNA\_N  
ClientSampleId :  
BHBX9

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_N\Methods\SFAM-EPA-BN052924.MA.M  
Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 26 9-Octadecenamide, (Z)- Concentration Rank 7

| R.T.      | EstConc                | Area    | Relative to ISTD | R.T.        |      |
|-----------|------------------------|---------|------------------|-------------|------|
| 22.668    | 4.50 ng/ul             | 1754620 | Chrysene-d12     | 21.304      |      |
| Hit# of 5 | Tentative ID           | MW      | MolForm          | CAS#        | Qual |
| 1         | 9-Octadecenamide, (Z)- | 281     | C18H35NO         | 000301-02-0 | 94   |
| 2         | 13-Docosenamide, (Z)-  | 337     | C22H43NO         | 000112-84-5 | 93   |
| 3         | 9-Octadecenamide, (Z)- | 281     | C18H35NO         | 000301-02-0 | 93   |
| 4         | 9-Octadecenamide, (Z)- | 281     | C18H35NO         | 000301-02-0 | 91   |
| 5         | 9-Octadecenamide, (Z)- | 281     | C18H35NO         | 000301-02-0 | 91   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN060824\  
 Data File : BN031856.D  
 Acq On : 08 Jun 2024 13:14  
 Operator : MA/JU  
 Sample : P2634-15  
 Misc :  
 ALS Vial : 7 Sample Multiplier: 1

Instrument :  
 BNA\_N  
 ClientSampleId :  
 BHBX9

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_N\Methods\SFAM-EPA-BN052924.MA.M  
 Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L  
 TIC Integration Parameters: LSCINT.P

| TIC Top Hit name   | RT     | EstConc | Units | Response | --Internal Standard-- |        |         |      |
|--------------------|--------|---------|-------|----------|-----------------------|--------|---------|------|
|                    |        |         |       |          | #                     | RT     | Resp    | Conc |
| unknown-01         | 3.693  | 3.3     | ng/ul | 394278   | 1                     | 7.669  | 2381800 | 20.0 |
| Styrene            | 5.675  | 5.6     | ng/ul | 669865   | 1                     | 7.669  | 2381800 | 20.0 |
| 1-Phenoxypropan... | 11.198 | 2.8     | ng/ul | 491336   | 2                     | 10.451 | 3554620 | 20.0 |
| (DEL) Alkane: S... | 16.034 | 2.5     | ng/ul | 659655   | 4                     | 17.063 | 5357360 | 20.0 |
| Benzene, 1,1'-(... | 16.669 | 3.6     | ng/ul | 963518   | 4                     | 17.063 | 5357360 | 20.0 |
| 9H-Fluoren-9-one   | 16.722 | 2.5     | ng/ul | 659185   | 4                     | 17.063 | 5357360 | 20.0 |
| 2-Methyldibenzo... | 17.645 | 2.4     | ng/ul | 642225   | 4                     | 17.063 | 5357360 | 20.0 |
| Phenanthrene, 2... | 17.939 | 4.7     | ng/ul | 1251600  | 4                     | 17.063 | 5357360 | 20.0 |
| n-Hexadecanoic ... | 17.969 | 4.4     | ng/ul | 1189820  | 4                     | 17.063 | 5357360 | 20.0 |
| 1H-Cyclopropa[1... | 18.128 | 5.9     | ng/ul | 1580260  | 4                     | 17.063 | 5357360 | 20.0 |
| Anthracene, 2-m... | 18.169 | 2.9     | ng/ul | 775932   | 4                     | 17.063 | 5357360 | 20.0 |
| Naphthalene, 2-... | 18.445 | 3.7     | ng/ul | 988100   | 4                     | 17.063 | 5357360 | 20.0 |
| 9,10-Anthracene... | 18.486 | 2.7     | ng/ul | 715658   | 4                     | 17.063 | 5357360 | 20.0 |
| Phenanthrene, 3... | 18.739 | 2.4     | ng/ul | 646847   | 4                     | 17.063 | 5357360 | 20.0 |
| Phenanthrene, 2... | 18.863 | 4.5     | ng/ul | 1192940  | 4                     | 17.063 | 5357360 | 20.0 |
| Cyclopenta(def)... | 19.004 | 3.2     | ng/ul | 860082   | 4                     | 17.063 | 5357360 | 20.0 |
| Pyrene, 1-methyl-  | 19.816 | 3.2     | ng/ul | 1264840  | 5                     | 21.304 | 7796100 | 20.0 |
| 11H-Benzo[b]flu... | 19.992 | 2.2     | ng/ul | 849534   | 5                     | 21.304 | 7796100 | 20.0 |
| 11H-Benzo[a]flu... | 20.739 | 2.4     | ng/ul | 945004   | 5                     | 21.304 | 7796100 | 20.0 |
| Cyclohexane, 1,... | 20.968 | 7.0     | ng/ul | 2714290  | 5                     | 21.304 | 7796100 | 20.0 |
| 1H-Indole, 1-me... | 21.504 | 7.0     | ng/ul | 2737540  | 5                     | 21.304 | 7796100 | 20.0 |
| unknown-02         | 21.569 | 8.2     | ng/ul | 3183340  | 5                     | 21.304 | 7796100 | 20.0 |
| unknown-03         | 21.621 | 2.5     | ng/ul | 972573   | 5                     | 21.304 | 7796100 | 20.0 |
| unknown-04         | 21.663 | 3.2     | ng/ul | 1252400  | 5                     | 21.304 | 7796100 | 20.0 |
| (DEL) Alkane: S... | 22.051 | 3.0     | ng/ul | 1166930  | 5                     | 21.304 | 7796100 | 20.0 |
| 9-Octadecenamid... | 22.668 | 4.5     | ng/ul | 1754620  | 5                     | 21.304 | 7796100 | 20.0 |